

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100424\
 Data File : BF139669.D
 Acq On : 04 Oct 2024 20:44
 Operator : RC/JU
 Sample : P4242-01MS
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-QMS

Quant Time: Oct 04 22:55:47 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 02 04:58:47 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	85047	20.000	ng	0.00
21) Naphthalene-d8	8.151	136	297861	20.000	ng	0.00
39) Acenaphthene-d10	9.904	164	139846	20.000	ng	0.00
64) Phenanthrene-d10	11.386	188	230700	20.000	ng	0.00
76) Chrysene-d12	14.027	240	166417	20.000	ng	0.00
86) Perylene-d12	15.498	264	135056	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.492	112	508275	103.586	ng	0.00
7) Phenol-d6	6.504	99	653844	99.088	ng	0.00
23) Nitrobenzene-d5	7.428	82	421848	71.706	ng	0.00
42) 2,4,6-Tribromophenol	10.692	330	125211	92.750	ng	0.00
45) 2-Fluorobiphenyl	9.222	172	701471	77.005	ng	0.00
79) Terphenyl-d14	12.968	244	621912	61.319	ng	-0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.716	88	93709	44.173	ng	99
3) Pyridine	3.463	79	222487	43.123	ng	98
4) n-Nitrosodimethylamine	3.422	42	156250	46.266	ng	98
6) Aniline	6.534	93	217201	37.094	ng	96
8) 2-Chlorophenol	6.651	128	240154	45.682	ng	99
9) Benzaldehyde	6.416	77	85364	24.422	ng	99
10) Phenol	6.516	94	315350	45.450	ng	99
11) bis(2-Chloroethyl)ether	6.604	93	252140	44.632	ng	99
12) 1,3-Dichlorobenzene	6.804	146	260435	44.037	ng	99
13) 1,4-Dichlorobenzene	6.886	146	263683	44.022	ng	99
14) 1,2-Dichlorobenzene	7.034	146	246368	43.887	ng	100
15) Benzyl Alcohol	7.010	79	212289	42.106	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.139	45	435242	44.503	ng	89
17) 2-Methylphenol	7.122	107	190624	43.435	ng	99
18) Hexachloroethane	7.375	117	89656	40.130	ng	98
19) n-Nitroso-di-n-propyla...	7.281	70	179553	42.324	ng	99
20) 3+4-Methylphenols	7.275	107	234943	42.620	ng	92
22) Acetophenone	7.275	105	327213	46.280	ng	97
24) Nitrobenzene	7.451	77	269158	44.183	ng	99
25) Isophorone	7.686	82	466052	44.603	ng	100
26) 2-Nitrophenol	7.763	139	120015	45.827	ng	98
27) 2,4-Dimethylphenol	7.804	122	172820	53.905	ng	99
28) bis(2-Chloroethoxy)met...	7.898	93	280624	44.975	ng	99
29) 2,4-Dichlorophenol	8.010	162	187285	44.786	ng	99
30) 1,2,4-Trichlorobenzene	8.086	180	203233	42.974	ng	100
31) Naphthalene	8.169	128	667152	43.807	ng	100
32) Benzoic acid	7.922	122	116489	36.540	ng	98
33) 4-Chloroaniline	8.233	127	49860	9.672	ng	99
34) Hexachlorobutadiene	8.286	225	132017	43.137	ng	100
35) Caprolactam	8.581	113	56629	41.287	ng	99
36) 4-Chloro-3-methylphenol	8.710	107	191024	40.353	ng	99
37) 2-Methylnaphthalene	8.863	142	412375	41.576	ng	100
38) 1-Methylnaphthalene	8.963	142	381977	39.398	ng	100
40) 1,2,4,5-Tetrachloroben...	9.028	216	195235	50.275	ng	98
41) Hexachlorocyclopentadiene	9.010	237	80655	67.736	ng	100
43) 2,4,6-Trichlorophenol	9.139	196	123389	47.133	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100424\
 Data File : BF139669.D
 Acq On : 04 Oct 2024 20:44
 Operator : RC/JU
 Sample : P4242-01MS
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-QMS

Quant Time: Oct 04 22:55:47 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 02 04:58:47 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.186	196	127079	45.013	ng	99
46) 1,1'-Biphenyl	9.322	154	499595	48.032	ng	99
47) 2-Chloronaphthalene	9.351	162	365257	46.842	ng	98
48) 2-Nitroaniline	9.445	65	132770	47.433	ng	98
49) Acenaphthylene	9.763	152	577558	48.705	ng	100
50) Dimethylphthalate	9.622	163	441217	48.197	ng	99
51) 2,6-Dinitrotoluene	9.686	165	89191	43.130	ng	98
52) Acenaphthene	9.933	154	362403	44.517	ng	99
53) 3-Nitroaniline	9.857	138	66857	31.637	ng	99
54) 2,4-Dinitrophenol	9.963	184	39657	35.654	ng	97
55) Dibenzofuran	10.110	168	510767	44.811	ng	99
56) 4-Nitrophenol	10.022	139	131788	81.777	ng	99
57) 2,4-Dinitrotoluene	10.092	165	113154	41.862	ng	98
58) Fluorene	10.451	166	384911	42.398	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.227	232	98486	43.192	ng	97
60) Diethylphthalate	10.322	149	428602	45.328	ng	100
61) 4-Chlorophenyl-phenyle...	10.439	204	192517	42.388	ng	99
62) 4-Nitroaniline	10.469	138	74824	34.615	ng	97
63) Azobenzene	10.604	77	443201	46.608	ng	95
65) 4,6-Dinitro-2-methylph...	10.498	198	28186	21.632	ng	99
66) n-Nitrosodiphenylamine	10.563	169	337487	49.629	ng	99
67) 4-Bromophenyl-phenylether	10.933	248	109282	46.201	ng	99
68) Hexachlorobenzene	10.998	284	116933	44.550	ng	98
69) Atrazine	11.086	200	101929	55.866	ng	99
70) Pentachlorophenol	11.192	266	131153	83.917	ng	100
71) Phenanthrene	11.416	178	517256	47.552	ng	99
72) Anthracene	11.463	178	522500	48.554	ng	100
73) Carbazole	11.621	167	488580	47.271	ng	100
74) Di-n-butylphthalate	11.945	149	607950	48.371	ng	99
75) Fluoranthene	12.598	202	572976	48.186	ng	100
77) Benzidine	12.727	184	302643	103.101	ng	100
78) Pyrene	12.827	202	591197	39.720	ng	99
80) Butylbenzylphthalate	13.445	149	254307	47.767	ng	99
81) Benzo(a)anthracene	14.015	228	524614	47.948	ng	100
82) 3,3'-Dichlorobenzidine	13.980	252	174376	52.076	ng	98
83) Chrysene	14.057	228	466472	46.633	ng	99
84) Bis(2-ethylhexyl)phtha...	14.004	149	356602	53.641	ng	99
85) Di-n-octyl phthalate	14.615	149	584252	59.821	ng	100
87) Indeno(1,2,3-cd)pyrene	16.968	276	338558	42.591	ng	99
88) Benzo(b)fluoranthene	15.068	252	382710	43.826	ng	99
89) Benzo(k)fluoranthene	15.098	252	378084	51.041	ng	99
90) Benzo(a)pyrene	15.433	252	340660	49.475	ng	100
91) Dibenzo(a,h)anthracene	16.986	278	280328	42.519	ng	98
92) Benzo(g,h,i)perylene	17.409	276	250172	37.835	ng	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

