

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100722\
 Data File : BF130582.D
 Acq On : 07 Oct 2022 19:08
 Operator : CG\JU
 Sample : N5018-02MS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SASP2-T-3-(5-10)MS

Quant Time: Oct 10 01:44:50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 07 02:06:46 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Christian Giraldo 10/10/2022
 Supervised By : Jagrut Upadhyay 10/10/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.616	152	82221	20.000	ng	0.00
21) Naphthalene-d8	7.898	136	308950	20.000	ng	# 0.00
39) Acenaphthene-d10	9.645	164	175152	20.000	ng	0.00
64) Phenanthrene-d10	11.122	188	256469	20.000	ng	-0.01
76) Chrysene-d12	13.757	240	172003	20.000	ng	0.00
86) Perylene-d12	15.139	264	239956	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.222	112	673157	142.003	ng	0.00
7) Phenol-d6	6.263	99	799514	134.794	ng	0.00
23) Nitrobenzene-d5	7.187	82	429242	70.980	ng	0.00
42) 2,4,6-Tribromophenol	10.433	330	164543	98.042	ng	0.00
45) 2-Fluorobiphenyl	8.975	172	845482	70.292	ng	0.00
79) Terphenyl-d14	12.710	244	758794	73.076	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.240	88	118055	54.226	ng	99
3) Pyridine	2.910	79	251772	44.332	ng	93
4) n-Nitrosodimethylamine	2.869	42	136362	44.147	ng	84
6) Aniline	6.281	93	341160	46.682	ng	# 46
8) 2-Chlorophenol	6.398	128	310135	57.433	ng	97
9) Benzaldehyde	6.163	77	143876	36.213	ng	94
10) Phenol	6.281	94	369627	53.176	ng	# 75
11) bis(2-Chloroethyl)ether	6.357	93	295886	59.016	ng	95
12) 1,3-Dichlorobenzene	6.551	146	282808	47.597	ng	98
13) 1,4-Dichlorobenzene	6.634	146	285422	47.106	ng	97
14) 1,2-Dichlorobenzene	6.781	146	268710	47.473	ng	99
15) Benzyl Alcohol	6.769	79	193327	41.109	ng	94
16) 2,2'-oxybis(1-Chloropr...	6.898	45	310183	42.418	ng	86
17) 2-Methylphenol	6.887	107	208983	47.608	ng	97
18) Hexachloroethane	7.122	117	109828	45.984	ng	97
19) n-Nitroso-di-n-propyla...	7.040	70	172267	43.041	ng	91
20) 3+4-Methylphenols	7.040	107	252506	44.280	ng	# 67
22) Acetophenone	7.034	105	358539	47.467	ng	# 91
24) Nitrobenzene	7.204	77	264883	43.137	ng	93
25) Isophorone	7.445	82	462490	47.450	ng	98
26) 2-Nitrophenol	7.522	139	134245	53.332	ng	# 83
27) 2,4-Dimethylphenol	7.563	122	207600	53.563	ng	98
28) bis(2-Chloroethoxy)met...	7.657	93	293283	53.437	ng	99
29) 2,4-Dichlorophenol	7.763	162	197313	46.457	ng	97
30) 1,2,4-Trichlorobenzene	7.839	180	220921	48.111	ng	97
31) Naphthalene	7.922	128	730954	45.924	ng	99
32) Benzoic acid	7.692	122	62892	20.983	ng	90
33) 4-Chloroaniline	7.975	127	179138	29.592	ng	97
34) Hexachlorobutadiene	8.034	225	143438	48.877	ng	99
35) Caprolactam	8.345	113	54908m	45.095	ng	
36) 4-Chloro-3-methylphenol	8.469	107	214332	45.528	ng	100
37) 2-Methylnaphthalene	8.610	142	479237	47.223	ng	100
38) 1-Methylnaphthalene	8.710	142	450678	46.228	ng	99
40) 1,2,4,5-Tetrachloroben...	8.775	216	217870	45.580	ng	99
41) Hexachlorocyclopentadiene	8.757	237	173225	67.688	ng	99
43) 2,4,6-Trichlorophenol	8.892	196	131700	39.478	ng	95

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44) 2,4,5-Trichlorophenol	8.934	196	138413	38.449	ng	97	10/10/2022
46) 1,1'-Biphenyl	9.075	154	643322	49.179	ng	99	Supervised By : Jagrut Upadhyay
47) 2-Chloronaphthalene	9.098	162	495034	46.782	ng	95	
48) 2-Nitroaniline	9.198	65	138548	39.256	ng	94	
49) Acenaphthylene	9.510	152	769134	48.477	ng	99	
50) Dimethylphthalate	9.381	163	544556	45.009	ng	99	
51) 2,6-Dinitrotoluene	9.445	165	118621	46.880	ng	# 76	10/10/2022
52) Acenaphthene	9.681	154	501634m	48.121	ng		
53) 3-Nitroaniline	9.610	138	92374	32.763	ng	95	
54) 2,4-Dinitrophenol	9.722	184	87754	65.189	ng	91	
55) Dibenzofuran	9.851	168	671673	46.324	ng	97	
56) 4-Nitrophenol	9.786	139	144504	70.367	ng	85	
57) 2,4-Dinitrotoluene	9.845	165	151374	46.810	ng	95	
58) Fluorene	10.192	166	492419	41.526	ng	100	
59) 2,3,4,6-Tetrachlorophenol	9.975	232	106647	37.849	ng	91	
60) Diethylphthalate	10.075	149	509321	41.979	ng	99	
61) 4-Chlorophenyl-phenyle...	10.186	204	231770	44.555	ng	99	
62) 4-Nitroaniline	10.222	138	110258m	38.928	ng		
63) Azobenzene	10.345	77	483474	38.699	ng	96	
65) 4,6-Dinitro-2-methylph...	10.251	198	66440	46.537	ng	97	
66) n-Nitrosodiphenylamine	10.310	169	396338	46.245	ng	99	
67) 4-Bromophenyl-phenylether	10.675	248	137325	48.537	ng	96	
68) Hexachlorobenzene	10.733	284	152378	47.512	ng	94	
69) Atrazine	10.839	200	123919	49.477	ng	96	
70) Pentachlorophenol	10.939	266	116752	67.830	ng	99	
71) Phenanthrene	11.151	178	773597	53.725	ng	100	
72) Anthracene	11.198	178	687757	49.495	ng	100	
73) Carbazole	11.363	167	596914	47.599	ng	99	
74) Di-n-butylphthalate	11.698	149	761788	50.374	ng	# 98	
75) Fluoranthene	12.333	202	988142	71.018	ng	98	
77) Benzidine	12.463	184	376599	75.424	ng	100	
78) Pyrene	12.563	202	927813	61.661	ng	99	
80) Butylbenzylphthalate	13.186	149	311076	55.806	ng	99	
81) Benzo(a)anthracene	13.745	228	682686	59.662	ng	99	
82) 3,3'-Dichlorobenzidine	13.716	252	149250	47.917	ng	99	
83) Chrysene	13.786	228	641289	59.025	ng	99	
84) Bis(2-ethylhexyl)phtha...	13.745	149	382514	59.909	ng	99	
85) Di-n-octyl phthalate	14.357	149	711974	72.904	ng	96	
87) Indeno(1,2,3-cd)pyrene	16.451	276	904902	53.179	ng	99	
88) Benzo(b)fluoranthene	14.757	252	910916	58.162	ng	100	
89) Benzo(k)fluoranthene	14.786	252	795296	51.621	ng	98	
90) Benzo(a)pyrene	15.086	252	807503	65.080	ng	99	
91) Dibenzo(a,h)anthracene	16.468	278	726283	52.028	ng	98	
92) Benzo(g,h,i)perylene	16.851	276	747896	53.186	ng	98	

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

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