

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091522\
 Data File : BF130271.D
 Acq On : 15 Sep 2022 15:06
 Operator : CG\JU
 Sample : N4613-02MS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SASP2-T-6-(18-23)MS

Quant Time: Sep 16 03:27:29 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 14 14:55:51 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Christian
 Giraldo

Supervised By : Jagrut
 Upadhyay
 09/16/2022

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.675	152	105115	20.000	ng	0.00
21) Naphthalene-d8	7.963	136	415922	20.000	ng	0.00
39) Acenaphthene-d10	9.710	164	216595	20.000	ng	0.00
64) Phenanthrene-d10	11.186	188	368669	20.000	ng	0.00
76) Chrysene-d12	13.821	240	186453	20.000	ng	# 0.00
86) Perylene-d12	15.209	264	173723	20.000	ng	0.00

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System Monitoring Compounds						
5) 2-Fluorophenol	5.287	112	807052	133.169	ng	0.01
7) Phenol-d6	6.316	99	1018895	134.367	ng	0.00
23) Nitrobenzene-d5	7.245	82	662354	81.358	ng	0.00
42) 2,4,6-Tribromophenol	10.498	330	323867	156.051	ng	0.00
45) 2-Fluorobiphenyl	9.039	172	1198253	80.560	ng	0.00
79) Terphenyl-d14	12.774	244	1053971	93.637	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	2.393	88	111336	40.001	ng	97
3) Pyridine	3.075	79	284601	39.198	ng	98
4) n-Nitrosodimethylamine	3.040	42	173893	44.036	ng	96
6) Aniline	6.339	93	407453	43.610	ng	95
8) 2-Chlorophenol	6.463	128	336453	48.736	ng	94
9) Benzaldehyde	6.228	77	211028	41.546	ng	96
10) Phenol	6.333	94	429775	48.363	ng	88
11) bis(2-Chloroethyl)ether	6.422	93	303756	47.390	ng	96
12) 1,3-Dichlorobenzene	6.616	146	359913	47.381	ng	99
13) 1,4-Dichlorobenzene	6.692	146	366585	47.324	ng	100
14) 1,2-Dichlorobenzene	6.845	146	347998	48.091	ng	98
15) Benzyl Alcohol	6.828	79	263876	43.890	ng	97
16) 2,2'-oxybis(1-Chloropr...	6.963	45	415160	44.409	ng	95
17) 2-Methylphenol	6.939	107	266174	47.430	ng	98
18) Hexachloroethane	7.186	117	143528	47.005	ng	98
19) n-Nitroso-di-n-propyla...	7.104	70	240423	46.987	ng	96
20) 3+4-Methylphenols	7.092	107	343571	47.128	ng	# 88
22) Acetophenone	7.098	105	484474	47.644	ng	# 95
24) Nitrobenzene	7.269	77	371689	44.963	ng	94
25) Isophorone	7.504	82	652258	49.708	ng	100
26) 2-Nitrophenol	7.580	139	175595	51.817	ng	94
27) 2,4-Dimethylphenol	7.622	122	298291	57.169	ng	99
28) bis(2-Chloroethoxy)met...	7.722	93	384521	52.042	ng	100
29) 2,4-Dichlorophenol	7.822	162	279882	48.949	ng	98
30) 1,2,4-Trichlorobenzene	7.904	180	283372	45.839	ng	95
31) Naphthalene	7.980	128	980283	45.748	ng	99
32) Benzoic acid	7.757	122	190288m	47.159	ng	
33) 4-Chloroaniline	8.033	127	233245	28.620	ng	99
34) Hexachlorobutadiene	8.098	225	181224	45.871	ng	100
35) Caprolactam	8.416	113	90396m	55.147	ng	
36) 4-Chloro-3-methylphenol	8.522	107	317243	50.056	ng	97
37) 2-Methylnaphthalene	8.675	142	639583	46.814	ng	99
38) 1-Methylnaphthalene	8.775	142	612234	46.648	ng	99
40) 1,2,4,5-Tetrachloroben...	8.839	216	288255	48.766	ng	98
41) Hexachlorocyclopentadiene	8.827	237	390352	123.346	ng	100
43) 2,4,6-Trichlorophenol	8.951	196	199642	48.394	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	8.992	196	214459	48.175	ng	99
46) 1,1'-Biphenyl	9.139	154	782182	48.354	ng	98
47) 2-Chloronaphthalene	9.163	162	613981	46.920	ng	98
48) 2-Nitroaniline	9.263	65	201485	46.165	ng	94
49) Acenaphthylene	9.574	152	927391	47.268	ng	100
50) Dimethylphthalate	9.445	163	729701	48.772	ng	100
51) 2,6-Dinitrotoluene	9.504	165	159710	51.042	ng	96
52) Acenaphthene	9.745	154	668648m	51.870	ng	09/16/2022
53) 3-Nitroaniline	9.669	138	117744	33.770	ng	100
54) 2,4-Dinitrophenol	9.780	184	150630	88.059	ng	# 82
55) Dibenzofuran	9.922	168	844127	47.078	ng	97
56) 4-Nitrophenol	9.839	139	247343	97.399	ng	94
57) 2,4-Dinitrotoluene	9.904	165	214462	53.630	ng	97
58) Fluorene	10.257	166	689556	47.025	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.033	232	163088	46.805	ng	98
60) Diethylphthalate	10.145	149	713814	47.577	ng	99
61) 4-Chlorophenyl-phenylether	10.257	204	318964	49.585	ng	95
62) 4-Nitroaniline	10.286	138	163886	46.790	ng	97
63) Azobenzene	10.416	77	686993	44.468	ng	96
65) 4,6-Dinitro-2-methylph...	10.310	198	92724	45.181	ng	97
66) n-Nitrosodiphenylamine	10.374	169	579273	47.020	ng	99
67) 4-Bromophenyl-phenylether	10.739	248	201244	49.482	ng	93
68) Hexachlorobenzene	10.798	284	223370	48.451	ng	# 89
69) Atrazine	10.904	200	194042	53.896	ng	99
70) Pentachlorophenol	10.998	266	232524	93.977	ng	99
71) Phenanthrene	11.216	178	940335	45.430	ng	100
72) Anthracene	11.268	178	942672	47.194	ng	99
73) Carbazole	11.427	167	828614	45.966	ng	100
74) Di-n-butylphthalate	11.763	149	1032559	47.500	ng	99
75) Fluoranthene	12.398	202	865565	43.276	ng	98
77) Benzidine	12.527	184	381036	70.398	ng	99
78) Pyrene	12.627	202	842520	51.653	ng	99
80) Butylbenzylphthalate	13.251	149	320076	52.970	ng	97
81) Benzo(a)anthracene	13.809	228	576749	46.498	ng	100
82) 3,3'-Dichlorobenzidine	13.780	252	164771	48.800	ng	98
83) Chrysene	13.845	228	532066	45.177	ng	99
84) Bis(2-ethylhexyl)phtha...	13.809	149	378813	54.731	ng	98
85) Di-n-octyl phthalate	14.421	149	528991	49.969	ng	98
87) Indeno(1,2,3-cd)pyrene	16.545	276	628539	51.020	ng	99
88) Benzo(b)fluoranthene	14.821	252	520810	45.932	ng	98
89) Benzo(k)fluoranthene	14.851	252	493165m	44.214	ng	
90) Benzo(a)pyrene	15.156	252	457541	50.934	ng	98
91) Dibenzo(a,h)anthracene	16.562	278	540420	53.474	ng	99
92) Benzo(g,h,i)perylene	16.950	276	543814	53.417	ng	98

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

