

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120919\
 Data File : BP001280.D
 Acq On : 09 Dec 2019 20:16
 Operator : JU
 Sample : K6177-08MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EP-6MSD

Manual Integrations
 APPROVED

mohammad
 12/10/2019 3:47:01 PM

Quant Time: Dec 10 05:30:21 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP112719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 05 12:34:47 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.31	152	63731	20.00	ng	0.00
21) Naphthalene-d8	10.35	136	236047	20.00	ng	0.00
39) Acenaphthene-d10	13.20	164	137781	20.00	ng	0.00
64) Phenanthrene-d10	15.60	188	270787	20.00	ng	0.00
76) Chrysene-d12	19.25	240	221082	20.00	ng	0.00
87) Perylene-d12	21.02	264	236961	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	6.22	112	391573	101.94	ng	0.01
7) Phenol-d6	7.76	99	544376	106.37	ng	0.00
23) Nitrobenzene-d5	9.19	82	373879	68.72	ng	0.00
42) 2,4,6-Tribromophenol	14.50	330	157034	118.91	ng	0.00
45) 2-Fluorobiphenyl	12.12	172	644163	68.43	ng	-0.01
79) Terphenyl-d14	17.89	244	812572	68.00	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.60	88	59141	32.961	ng	97
3) Pyridine	3.45	79	146037	29.331	ng	# 90
4) n-Nitrosodimethylamine	3.35	42	102461	47.557	ng	86
6) Aniline	7.78	93	73551	10.840	ng	# 1
8) 2-Chlorophenol	7.97	128	158554	39.894	ng	97
9) Benzaldehyde	7.60	77	70121	19.945	ng	98
10) Phenol	7.78	94	223819	38.449	ng	90
11) bis(2-Chloroethyl)ether	7.91	93	164493	36.288	ng	99
12) 1,3-Dichlorobenzene	8.22	146	179766	38.161	ng	97
13) 1,4-Dichlorobenzene	8.33	146	184068	38.789	ng	97
14) 1,2-Dichlorobenzene	8.57	146	172750	38.681	ng	99
15) Benzyl Alcohol	8.54	79	185292	44.107	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.77	45	254222	34.884	ng	100
17) 2-Methylphenol	8.73	107	139580	38.286	ng	97
18) Hexachloroethane	9.11	117	77576	39.518	ng	98
19) n-Nitroso-di-n-propylamine	8.97	70	143052	38.738	ng	100
20) 3+4-Methylphenols	8.97	107	182841	39.786	ng	98
22) Acetophenone	8.96	105	267481	38.555	ng	# 98
24) Nitrobenzene	9.22	77	215340	39.803	ng	99
25) Isophorone	9.62	82	369432	37.866	ng	98
26) 2-Nitrophenol	9.73	139	80586	40.666	ng	96
27) 2,4-Dimethylphenol	9.83	122	140908	44.891	ng	98
28) bis(2-Chloroethoxy)methane	9.98	93	209356	34.202	ng	100
29) 2,4-Dichlorophenol	10.12	162	143215	40.088	ng	99
30) 1,2,4-Trichlorobenzene	10.26	180	159793	38.295	ng	97
31) Naphthalene	10.37	128	465695	38.629	ng	99
32) Benzoic acid	10.00	122	86298	38.029	ng	95
33) 4-Chloroaniline	10.47	127	28612	5.469	ng	97
34) Hexachlorobutadiene	10.60	225	109106	41.566	ng	98
35) Caprolactam	11.04	113	45463m	36.924	ng	
36) 4-Chloro-3-methylphenol	11.29	107	172544	40.736	ng	99
37) 2-Methylnaphthalene	11.51	142	321671	39.103	ng	98
38) 1-Methylnaphthalene	11.67	142	296417	38.145	ng	99
40) 1,2,4,5-Tetrachlorobenzene	11.79	216	173161	39.880	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	11.78	237	214949	107.307	nq	97
43) 2,4,6-Trichlorophenol	11.97	196	109896	38.763	nq	96
44) 2,4,5-Trichlorophenol	12.03	196	118093	40.202	nq	99
46) 1,1'-Biphenyl	12.28	154	407267	38.247	nq	99
47) 2-Chloronaphthalene	12.30	162	318394	37.434	nq	99
48) 2-Nitroaniline	12.47	65	126951	38.083	nq	96
49) Acenaphthylene	12.97	152	503000	39.453	nq	100
50) Dimethylphthalate	12.80	163	438846	42.587	nq	98
51) 2,6-Dinitrotoluene	12.88	165	82928	37.594	nq	96
52) Acenaphthene	13.26	154	289989	37.812	nq	98
53) 3-Nitroaniline	13.13	138	23310	9.407	nq	98
54) 2,4-Dinitrophenol	13.32	184	70604	78.506	nq	86
55) Dibenzofuran	13.55	168	455310	39.508	nq	100
56) 4-Nitrophenol	13.45	139	137345	78.220	nq	94
57) 2,4-Dinitrotoluene	13.54	165	114353	41.358	nq	91
58) Fluorene	14.11	166	361910	39.191	nq	100
59) 2,3,4,6-Tetrachlorophenol	13.75	232	97503	40.379	nq	97
60) Diethylphthalate	13.97	149	400386	37.525	nq	99
61) 4-Chlorophenyl-phenylether	14.13	204	183379	38.997	nq	100
62) 4-Nitroaniline	12.47	138	100076	34.834	nq	98
63) Azobenzene	14.39	77	436989	37.889	nq	96
65) 4,6-Dinitro-2-methylphenol	14.20	198	48589	42.644	nq	90
66) n-Nitrosodiphenylamine	14.32	169	316410	38.457	nq	99
67) 4-Bromophenyl-phenylether	14.92	248	110040	37.431	nq	98
68) Hexachlorobenzene	15.00	284	119802	37.067	nq	99
69) Atrazine	15.21	200	105080	41.828	nq	96
70) Pentachlorophenol	15.32	266	144366	85.702	nq	99
71) Phenanthrene	15.64	178	549504	38.599	nq	100
72) Anthracene	15.72	178	559772	39.893	nq	99
73) Carbazole	15.96	167	501550	39.281	nq	99
74) Di-n-butylphthalate	16.52	149	655598	38.189	nq	100
75) Fluoranthene	17.35	202	624103	41.410	nq	99
77) Benzidine	17.54	184	179241	31.401	nq	98
78) Pyrene	17.66	202	631250	36.252	nq	100
80) Butylbenzylphthalate	18.54	149	288651	35.890	nq	99
81) Benzo(a)anthracene	19.23	228	560683	36.578	nq	99
82) 3,3'-Dichlorobenzidine	19.20	252	112714	22.258	nq	98
83) Chrysene	19.28	228	548724	39.977	nq	99
84) Bis(2-ethylhexyl)phthalate	19.29	149	411475	37.211	nq	100
85) Di-n-octyl phthalate	20.07	149	707499	36.018	nq	100
86) Indeno(1,2,3-cd)pyrene	22.66	276	595189	36.793	nq	98
88) Benzo(b)fluoranthene	20.53	252	550894	38.062	nq	99
89) Benzo(k)fluoranthene	20.57	252	534784	38.363	nq	99
90) Benzo(a)pyrene	20.95	252	528413	39.636	nq	98
91) Dibenzo(a,h)anthracene	22.69	278	499359	38.275	nq	98
92) Benzo(g,h,i)perylene	23.16	276	485025	37.650	nq	97

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

