

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112520\
 Data File : BP004108.D
 Acq On : 25 Nov 2020 21:21
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

mohammad
 11/27/2020 11:45:22 AM

Quant Time: Nov 26 00:32:43 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 23 16:42:41 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.263	152	88531	20.00	ng	0.00
21) Naphthalene-d8	11.099	136	348399	20.00	ng	# 0.00
39) Acenaphthene-d10	14.887	164	227444	20.00	ng	0.00
64) Phenanthrene-d10	17.622	188	511022	20.00	ng	# 0.00
76) Chrysene-d12	21.657	240	545592	20.00	ng	0.00
86) Perylene-d12	24.127	264	630703	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.799	112	435597	82.12	ng	0.00
7) Phenol-d6	7.452	99	637773	83.76	ng	0.00
23) Nitrobenzene-d5	9.428	82	567148	79.73	ng	0.00
42) 2,4,6-Tribromophenol	16.375	330	252334	88.70	ng	0.00
45) 2-Fluorobiphenyl	13.510	172	1305649	80.25	ng	0.00
79) Terphenyl-d14	20.122	244	2216525	78.49	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.475	88	85667	37.43	ng	# 60
3) Pyridine	3.917	79	268837	40.05	ng	# 63
4) n-Nitrosodimethylamine	3.817	42	126703	41.00	ng	# 53
6) Aniline	7.569	93	364727	42.18	ng	# 83
8) 2-Chlorophenol	7.840	128	234500	41.64	ng	# 80
9) Benzaldehyde	7.369	77	136637	39.96	ng	# 82
10) Phenol	7.475	94	303676	41.33	ng	83
11) bis(2-Chloroethyl)ether	7.657	93	235427	39.86	ng	# 80
12) 1,3-Dichlorobenzene	8.157	146	278784	40.39	ng	# 94
13) 1,4-Dichlorobenzene	8.299	146	281524	40.46	ng	94
14) 1,2-Dichlorobenzene	8.634	146	270712	40.73	ng	97
15) Benzyl Alcohol	8.505	79	226328	42.91	ng	# 87
16) 2,2'-oxybis(1-Chloropr...	8.793	45	382347	38.44	ng	83
17) 2-Methylphenol	8.740	107	206306	42.76	ng	# 94
18) Hexachloroethane	9.375	117	104311	42.18	ng	96
19) n-Nitroso-di-n-propyla...	9.069	70	194527	42.74	ng	# 78
20) 3+4-Methylphenols	9.063	107	280121	43.50	ng	91
22) Acetophenone	9.087	105	373905	39.98	ng	# 84
24) Nitrobenzene	9.475	77	289616	40.97	ng	# 83
25) Isophorone	9.999	82	514551	42.58	ng	# 91
26) 2-Nitrophenol	10.199	139	133359	48.31	ng	# 77
27) 2,4-Dimethylphenol	10.269	122	190003	41.27	ng	91
28) bis(2-Chloroethoxy)met...	10.469	93	329523	39.74	ng	97
29) 2,4-Dichlorophenol	10.763	162	235211	42.39	ng	95
30) 1,2,4-Trichlorobenzene	10.963	180	274466	40.60	ng	98
31) Naphthalene	11.146	128	752724	40.26	ng	100
32) Benzoic acid	10.451	122	87015	30.57	ng	# 82
33) 4-Chloroaniline	11.240	127	331675	41.74	ng	# 90
34) Hexachlorobutadiene	11.451	225	176650	40.07	ng	98
35) Caprolactam	11.981	113	86439	50.38	ng	# 16
36) 4-Chloro-3-methylphenol	12.381	107	260634	43.51	ng	90
37) 2-Methylnaphthalene	12.734	142	546142	40.82	ng	98
38) 1-Methylnaphthalene	12.951	142	522136	40.92	ng	98
40) 1,2,4,5-Tetrachloroben...	13.116	216	302666	40.21	ng	99
41) Hexachlorocyclopentadiene	13.104	237	113668	29.93	ng	99
43) 2,4,6-Trichlorophenol	13.345	196	196157	42.15	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	13.434	196	203167	41.47	ng	#	95
46) 1,1'-Biphenyl	13.722	154	700382	39.67	ng		98
47) 2-Chloronaphthalene	13.769	162	587518	40.56	ng		99
48) 2-Nitroaniline	13.957	65	202525	46.85	ng	#	61
49) Acenaphthylene	14.610	152	886586	41.68	ng		100
50) Dimethylphthalate	14.316	163	747279	42.03	ng		100
51) 2,6-Dinitrotoluene	14.440	165	161908	44.38	ng	#	75
52) Acenaphthene	14.951	154	570928m	40.29	ng		
53) 3-Nitroaniline	14.775	138	179053	44.34	ng	#	75
54) 2,4-Dinitrophenol	14.992	184	49022	33.15	ng	#	83
55) Dibenzofuran	15.281	168	860428	40.87	ng		100
56) 4-Nitrophenol	15.122	139	121124	41.62	ng	#	65
57) 2,4-Dinitrotoluene	15.234	165	232988	47.57	ng	#	76
58) Fluorene	15.928	166	703025	41.70	ng		99
59) 2,3,4,6-Tetrachlorophenol	15.528	232	176105	39.87	ng		97
60) Diethylphthalate	15.669	149	743695	42.72	ng		99
61) 4-Chlorophenyl-phenyle...	15.904	204	376670	40.88	ng		98
62) 4-Nitroaniline	15.934	138	201402	47.84	ng		88
63) Azobenzene	16.198	77	665109	40.47	ng		85
65) 4,6-Dinitro-2-methylph...	16.016	198	99515	37.73	ng		86
66) n-Nitrosodiphenylamine	16.116	169	621911	39.75	ng		98
67) 4-Bromophenyl-phenylether	16.798	248	238694	39.92	ng		98
68) Hexachlorobenzene	16.963	284	272766	39.98	ng		98
69) Atrazine	17.051	200	208694	40.64	ng		99
70) Pentachlorophenol	17.298	266	133455	40.89	ng		98
71) Phenanthrene	17.663	178	1157482	40.36	ng		99
72) Anthracene	17.751	178	1136808	41.04	ng		99
73) Carbazole	18.010	167	1090808	42.45	ng		98
74) Di-n-butylphthalate	18.522	149	1308584	43.99	ng		99
75) Fluoranthene	19.598	202	1419541	42.90	ng		99
77) Benzidine	19.745	184	452247	35.14	ng		100
78) Pyrene	19.951	202	1462860	39.51	ng		99
80) Butylbenzylphthalate	20.763	149	632258	45.83	ng	#	86
81) Benzo(a)anthracene	21.639	228	1471985	40.91	ng		98
82) 3,3'-Dichlorobenzidine	21.551	252	534150	43.04	ng		99
83) Chrysene	21.698	228	1410626	40.81	ng		98
84) Bis(2-ethylhexyl)phtha...	21.510	149	871510	42.86	ng		99
85) Di-n-octyl phthalate	22.439	149	1509716	44.56	ng	#	87
87) Indeno(1,2,3-cd)pyrene	26.698	276	1945808	42.77	ng		98
88) Benzo(b)fluoranthene	23.380	252	1583559	38.37	ng		99
89) Benzo(k)fluoranthene	23.427	252	1472170	36.12	ng		98
90) Benzo(a)pyrene	24.021	252	1488943	38.97	ng		98
91) Dibenzo(a,h)anthracene	26.692	278	1623736	42.81	ng		98
92) Benzo(g,h,i)perylene	27.486	276	1604920	43.72	ng		97

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

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