

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN072618\
 Data File : BN002173.D
 Acq On : 26 Jul 2018 15:32
 Operator : JU/SJ
 Sample : J4086-01DL 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 1524DL

Quant Time: Jul 27 01:52:05 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-BN072518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 26 18:27:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	223386	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	982028	20.00	ng	0.00
38) Acenaphthene-d10	14.32	164	548041	20.00	ng	0.00
63) Phenanthrene-d10	17.07	188	1074876	20.00	ng	0.00
75) Chrysene-d12	21.27	240	839689	20.00	ng	0.00
86) Perylene-d12	23.49	264	817786	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.30	112	249004	17.95	ng	0.00
7) Phenol-d6	6.87	99	367621	18.88	ng	0.00
23) Nitrobenzene-d5	8.84	82	244433	12.76	ng	0.00
41) 2,4,6-Tribromophenol	15.82	330	150486	21.15	ng	0.00
44) 2-Fluorobiphenyl	12.94	172	606828	14.42	ng	0.00
78) Terphenyl-d14	19.70	244	777254	19.31	ng	0.00

Target Compounds

						Qvalue
8) 2-Chlorophenol	7.26	128	553910	34.421	ng	93
10) Phenol	6.90	94	581854	29.152	ng	98
13) 1,4-Dichlorobenzene	7.73	146	154802	8.251	ng	97
14) 1,2-Dichlorobenzene	8.03	146	147235	8.135	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.22	45	494420	18.847	ng	94
17) 2-Methylphenol	8.13	107	533927	39.184	ng	96
18) Hexachloroethane	8.76	117	111484	16.236	ng	87
19) n-Nitroso-di-n-propylamine	8.49	70	337541	23.899	ng	# 97
25) Isophorone	9.40	82	1167302	35.502	ng	96
26) 2-Nitrophenol	9.59	139	76798	8.011	ng	96
29) 2,4-Dichlorophenol	10.12	162	389971	25.083	ng	95
31) Naphthalene	10.51	128	436424	8.806	ng	97
34) Hexachlorobutadiene	10.80	225	153615	14.247	ng	97
42) 2,4,6-Trichlorophenol	12.75	196	360863	29.187	ng	99
43) 2,4,5-Trichlorophenol	12.82	196	271054	20.696	ng	97
46) 2-Chloronaphthalene	13.19	162	1339939	35.996	ng	99
48) Acenaphthylene	14.04	152	1943673	34.409	ng	98
49) Dimethylphthalate	13.78	163	349208	7.671	ng	100
50) 2,6-Dinitrotoluene	13.90	165	329213	32.551	ng	90
51) Acenaphthene	14.39	154	1262378	37.101	ng	99
54) Dibenzofuran	14.72	168	2286268	41.443	ng	96
57) Fluorene	15.37	166	1817428	43.709	ng	95
59) Diethylphthalate	15.15	149	1871635	40.393	ng	98
66) 4-Bromophenyl-phenylether	16.26	248	311912	24.788	ng	# 90
69) Pentachlorophenol	16.72	266	127588	14.241	ng	99
70) Phenanthrene	17.11	178	2015903	34.359	ng	100
71) Anthracene	17.20	178	671739	11.567	ng	97
79) Butylbenzylphthalate	20.41	149	682165	27.063	ng	92
80) Benzo(a)anthracene	21.25	228	1435277	28.044	ng	98
83) Bis(2-ethylhexyl)phthalate	21.19	149	791305	22.512	ng	96
84) Di-n-octyl phthalate	22.07	149	891374	15.993	ng	98
85) Indeno(1,2,3-cd)pyrene	25.72	276	569543	11.988	ng	# 93
87) Benzo(b)fluoranthene	22.83	252	1080717	22.392	ng	98
88) Benzo(k)fluoranthene	22.87	252	2152602	45.287	ng	# 98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
89) Benzo(a)pyrene	23.40	252	2112822	46.874	ng	97
90) Dibenzo(a,h)anthracene	25.73	278	616810	14.575	ng #	95

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

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