

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN052919\
 Data File : BN005927.D
 Acq On : 29 May 2019 13:18
 Operator : JU/SJ
 Sample : K2928-07
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A42C3

Quant Time: May 29 13:59:54 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN052819MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 29 02:28:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	160046	20.00	ng/ul	0.00
18) Naphthalene-d8	10.51	136	750894	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.37	164	456750	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.12	188	1107804	20.00	ng/ul	0.00
77) Chrysene-d12	21.33	240	1187155	20.00	ng/ul	0.00
85) Perylene-d12	23.60	264	1328291	20.00	ng/ul	0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.26	96	21800	5.99	ng/uL	0.00
5) Phenol-d5	6.89	99	388201	27.37	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.06	67	283471	35.03	ng/ul	0.00
9) 2-Chlorophenol-d4	7.26	132	335231	31.02	ng/ul	0.00
13) 4-Methylphenol-d8	8.42	113	230581	20.14	ng/ul	0.00
19) Nitrobenzene-d5	8.87	128	173664	36.03	ng/ul	0.00
22) 2-Nitrophenol-d4	9.59	143	170273	37.94	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.13	165	318472	30.05	ng/ul	0.00
29) 4-Chloroaniline-d4	10.65	131	385356	27.73	ng/ul	0.00
43) Dimethylphthalate-d6	13.77	166	1238021	36.77	ng/ul	0.00
46) Acenaphthylene-d8	14.06	160	1433996	34.05	ng/ul	0.00
51) 4-Nitrophenol-d4	14.56	143	128073	20.90	ng/ul	0.00
57) Fluorene-d10	15.36	176	1063636	37.67	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.48	200	90673	20.08	ng/ul	0.00
70) Anthracene-d10	17.22	188	1581667	33.55	ng/ul	0.00
78) Pyrene-d10	19.52	212	1899426	31.68	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.46	264	1932987	32.61	ng/ul	0.00

Target Compounds

					Qvalue	
6) Phenol	6.92	94	22831	1.562	ng/ul	97
28) Naphthalene	10.56	128	40791	1.145	ng/ul#	93
34) 2-Methylnaphthalene	12.17	142	29583	1.127	ng/ul	100
44) Dimethylphthalate	13.82	163	452278	13.439	ng/ul	99
69) Phenanthrene	17.16	178	402482	7.333	ng/ul	100
76) Fluoranthene	19.19	202	548431	9.360	ng/ul#	93
79) Pyrene	19.55	202	500656	6.658	ng/ul#	92
82) Benzo(a)anthracene	21.32	228	341636	4.578	ng/ul#	89
84) Chrysene	21.37	228	538434	7.736	ng/ul#	94
87) Benzo(b)fluoranthene	22.92	252	503518	6.749	ng/ul#	96
88) Benzo(k)fluoranthene	22.96	252	161183	2.350	ng/ul#	91
90) Benzo(a)pyrene	23.50	252	318460	4.526	ng/ul#	92
91) Indeno(1,2,3-cd)pyrene	25.87	276	182653	2.212	ng/ul#	89
92) Dibenzo(a,h)anthracene	25.87	278	75791	1.103	ng/ul#	65
93) Benzo(g,h,i)perylene	26.57	276	141083	2.029	ng/ul#	90

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

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