

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP022920\
 Data File : BP001965.D
 Acq On : 29 Feb 2020 18:30
 Operator : CG/JU
 Sample : L1615-14
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBDN8

Quant Time: Mar 02 02:43:04 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP022720MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Feb 27 14:59:37 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	296516	20.00	ng/ul	0.00
18) Naphthalene-d8	10.41	136	1189647	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.27	164	751210	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.03	188	1642011	20.00	ng/ul	0.00
78) Chrysene-d12	21.12	240	1637373	20.00	ng/ul	0.00
86) Perylene-d12	23.33	264	1857128	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.17	96	30075	4.36	ng/uL	0.00
5) Phenol-d5	6.82	99	536006	23.70	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.98	67	315572	26.35	ng/ul	0.00
9) 2-Chlorophenol-d4	7.18	132	479979	24.53	ng/ul	0.00
13) 4-Methylphenol-d8	8.35	113	448587	24.13	ng/ul	0.00
19) Nitrobenzene-d5	8.78	128	219085	23.97	ng/ul	0.00
22) 2-Nitrophenol-d4	9.50	143	238376	23.18	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.04	165	468783	23.67	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	622138	28.66	ng/ul	0.00
44) Dimethylphthalate-d6	13.69	166	1510805	26.57	ng/ul	0.00
47) Acenaphthylene-d8	13.96	160	1796559	26.42	ng/ul	0.00
52) 4-Nitrophenol-d4	14.47	143	189305	17.42	ng/ul	0.00
58) Fluorene-d10	15.27	176	1357001	27.59	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.39	200	154709	17.09	ng/ul	0.00
71) Anthracene-d10	17.13	188	2032011	27.53	ng/ul	0.00
79) Pyrene-d10	19.37	212	2373985	27.70	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.19	264	2595817	27.32	ng/ul	0.00

Target Compounds

				Ovalue	
6) Phenol	6.85	94	31032	1.322	ng/ul 98
45) Dimethylphthalate	13.74	163	287468	4.899	ng/ul 100

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

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