

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP021721\
 Data File : BP004800.D
 Acq On : 18 Feb 2021 06:55
 Operator : CG/JU
 Sample : M1407-19
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBGM7

Manual Integrations
 APPROVED

mohammad
 2/19/2021 12:52:32 PM

Quant Time: Feb 18 07:50:00 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP020821MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Feb 18 03:00:36 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	53269	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	194779	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.42	164	108907	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.17	188	215896	20.00	ng/ul	0.00
78) Chrysene-d12	21.26	240	236181	20.00	ng/ul	0.00
86) Perylene-d12	23.56	264	263919	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.26	96	6624	4.81	ng/uL	0.00
5) Phenol-d5	6.95	99	28903	6.76	ng/ul	0.02
7) Bis-(2-Chloroethyl)ether-d	7.10	67	67858	30.56	ng/ul	0.00
9) 2-Chlorophenol-d4	7.30	132	88211	27.23	ng/ul	0.00
13) 4-Methylphenol-d8	8.49	113	51646	15.35	ng/ul	0.02
19) Nitrobenzene-d5	8.93	128	49162	36.56	ng/ul	0.00
22) 2-Nitrophenol-d4	9.65	143	55833	37.83	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.19	165	88128	29.67	ng/ul	0.01
29) 4-Chloroaniline-d4	10.71	131	6647	2.00	ng/ul	0.01
44) Dimethylphthalate-d6	13.83	166	284947	36.04	ng/ul	0.00
47) Acenaphthylene-d8	14.11	160	360997	37.54	ng/ul	0.00
52) 4-Nitrophenol-d4	14.65	143	8774m	6.35	ng/ul	0.04
58) Fluorene-d10	15.41	176	249548	36.84	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.53	200	47586	36.46	ng/ul	0.00
71) Anthracene-d10	17.27	188	419264	44.04	ng/ul	0.00
79) Pyrene-d10	19.51	212	472868	42.77	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.42	264	584500	43.78	ng/ul	0.00

Target Compounds

				Ovalue
6) Phenol	6.98	94	6598	1.480 ng/ul 97
45) Dimethylphthalate	13.87	163	19801	2.429 ng/ul 97

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

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