

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP032921\
 Data File : BP005097.D
 Acq On : 30 Mar 2021 10:15
 Operator : CG/JU
 Sample : M1783-05DL 5X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0G11DL

Quant Time: Mar 30 10:45:07 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP032921MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Mar 29 15:02:46 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	30982	20.00	ng/ul	0.00
18) Naphthalene-d8	10.50	136	123355	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.36	164	81522	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.12	188	181945	20.00	ng/ul	0.00
78) Chrysene-d12	21.22	240	195073	20.00	ng/ul	0.00
86) Perylene-d12	23.50	264	194278	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
5) Phenol-d5	6.89	99	2536	0.95	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.05	67	6888	4.93	ng/ul	0.00
9) 2-Chlorophenol-d4	7.25	132	7174	3.82	ng/ul	0.00
13) 4-Methylphenol-d8	8.43	113	4218	2.07	ng/ul	0.00
19) Nitrobenzene-d5	8.88	128	4197	4.61	ng/ul	0.00
22) 2-Nitrophenol-d4	9.59	143	4548	4.52	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.13	165	7328	3.73	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
44) Dimethylphthalate-d6	13.78	166	30259	5.11	ng/ul	0.00
47) Acenaphthylene-d8	14.06	160	33206	4.53	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
58) Fluorene-d10	15.36	176	26270	5.13	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.49	200	4261	3.55	ng/ul	0.00
71) Anthracene-d10	17.22	188	47808	5.84	ng/ul	0.00
79) Pyrene-d10	19.47	212	56236	6.03	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.35	264	60131	6.03	ng/ul	-0.01

Target Compounds				Ovalue		
73) 1,2,3,4-Tetrachlorobenzene	13.16	216	4267	1.541	ng/uL	95

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

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