

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102921\
 Data File : BP007783.D
 Acq On : 29 Oct 2021 14:10
 Operator : CG/JU
 Sample : M4282-18DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFY99DL

Quant Time: Oct 29 14:39:24 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP102521.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 27 15:37:43 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	191246	20.000	ng/u1	0.00
20) Naphthalene-d8	10.734	136	756154	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.569	164	503158	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.322	188	1145466	20.000	ng/u1	0.00
79) Chrysene-d12	21.416	240	1350872	20.000	ng/u1	0.00
88) Perylene-d12	23.857	264	1451678	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	1579	0.375	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/u1	
7) Phenol-d5	7.093	99	14467	0.996	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.258	67	43881	5.441	ng/u1	0.00
11) 2-Chlorophenol-d4	7.458	132	46173	4.087	ng/u1	0.00
15) 4-Methylphenol-d8	8.640	113	27317	2.402	ng/u1	0.00
21) Nitrobenzene-d5	9.087	128	27593	5.355	ng/u1	0.00
24) 2-Nitrophenol-d4	9.810	143	25663	4.560	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.358	165	50940	4.447	ng/u1	0.00
31) 4-Chloroaniline-d4	10.869	131	62237	4.109	ng/u1	0.00
46) Dimethylphthalate-d6	13.975	166	203789	5.979	ng/u1	0.00
49) Acenaphthylene-d8	14.257	160	218441	5.229	ng/u1	0.00
54) 4-Nitrophenol-d4	14.769	143	5108	0.818	ng/u1	0.00
60) Fluorene-d10	15.563	176	189057	6.580	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	11811	1.928	ng/u1	0.00
73) Anthracene-d10	17.422	188	343814	7.304	ng/u1	0.00
81) Pyrene-d10	19.657	212	435095	6.918	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.698	264	468836	6.574	ng/u1	-0.01
Target Compounds						
2) 1,4-Dioxane	3.399	88	47948	10.187	ng/uL	Qvalue 93

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

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