

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102121\
 Data File : BP007569.D
 Acq On : 22 Oct 2021 00:53
 Operator : CG/JU
 Sample : M4217-10
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JDGG0

Quant Time: Oct 22 02:01:33 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 20 10:49:33 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	208093	20.000	ng/u1	0.00
20) Naphthalene-d8	10.752	136	859155	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.581	164	578258	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.334	188	1238946	20.000	ng/u1	0.00
79) Chrysene-d12	21.422	240	1266745	20.000	ng/u1	0.00
88) Perylene-d12	23.869	264	1296340	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.376	96	7911	1.501	ng/uL	0.00
4) Pyridine-d5	3.805	84	28246	1.958	ng/u1	0.01
7) Phenol-d5	7.105	99	139125	8.561	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	87245	9.001	ng/u1	0.00
11) 2-Chlorophenol-d4	7.475	132	115685	9.049	ng/u1	0.00
15) 4-Methylphenol-d8	8.652	113	109344	8.735	ng/u1	0.00
21) Nitrobenzene-d5	9.105	128	51960	9.557	ng/u1	0.00
24) 2-Nitrophenol-d4	9.828	143	49517	9.750	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	108672	8.773	ng/u1	0.00
31) 4-Chloroaniline-d4	10.881	131	139779	8.344	ng/u1	0.00
46) Dimethylphthalate-d6	13.993	166	352067	8.800	ng/u1	0.00
49) Acenaphthylene-d8	14.275	160	420601	8.366	ng/u1	0.00
54) 4-Nitrophenol-d4	14.763	143	21733	3.386	ng/u1	-0.01
60) Fluorene-d10	15.575	176	316747	9.461	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	15076	3.037	ng/u1	0.00
73) Anthracene-d10	17.434	188	501028	9.522	ng/u1	0.00
81) Pyrene-d10	19.663	212	607070	9.586	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.704	264	565003	8.759	ng/u1	-0.01

Target Compounds	Qvalue

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

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