

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122321\
 Data File : BP008542.D
 Acq On : 24 Dec 2021 01:01
 Operator : CG/JU
 Sample : M5126-17
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J19

Quant Time: Dec 24 04:15:48 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP122321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 23 14:48:11 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	557522	20.000	ng/u1	0.00
20) Naphthalene-d8	10.704	136	2363206	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.534	164	1688760	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.281	188	3682785	20.000	ng/u1	0.00
79) Chrysene-d12	21.363	240	3315252	20.000	ng/u1	0.00
88) Perylene-d12	23.792	264	2771860	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	56852	4.596	ng/uL	0.00
4) Pyridine-d5	3.758	84	174454	5.040	ng/u1	0.00
7) Phenol-d5	7.058	99	234959	5.283	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	848720	33.597	ng/u1	0.00
11) 2-Chlorophenol-d4	7.434	132	880950	24.899	ng/u1	0.00
15) 4-Methylphenol-d8	8.605	113	499313	13.994	ng/u1	0.00
21) Nitrobenzene-d5	9.057	128	615260	36.793	ng/u1	0.00
24) 2-Nitrophenol-d4	9.781	143	597676	38.980	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	1134101	30.132	ng/u1	0.00
31) 4-Chloroaniline-d4	10.834	131	1287885	24.789	ng/u1	0.00
46) Dimethylphthalate-d6	13.945	166	4723310	37.855	ng/u1	0.00
49) Acenaphthylene-d8	14.228	160	5616546	35.999	ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	102425	5.178	ng/u1	0.00
60) Fluorene-d10	15.528	176	4224668	38.997	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	518860	32.005	ng/u1	0.00
73) Anthracene-d10	17.381	188	7250721	41.547	ng/u1	0.00
81) Pyrene-d10	19.610	212	8180201	41.178	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.633	264	6101527	42.769	ng/u1	0.00

Target Compounds Qvalue

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

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