

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092923\
 Data File : BP017453.D
 Acq On : 29 Sep 2023 19:35
 Operator : MA/JU
 Sample : 04497-13
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBBP6

Quant Time: Sep 30 00:29:58 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 29 22:31:03 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	605183	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.793	136	2504581	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.645	164	1612976	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.434	188	3636590	20.000	ng/u1	-0.01
79) Chrysene-d12	21.563	240	3517533	20.000	ng/u1	-0.01
88) Perylene-d12	24.145	264	4022105	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	14887	1.010	ng/uL	0.00
4) Pyridine-d5	3.734	84	183104	4.445	ng/u1	0.00
7) Phenol-d5	7.111	99	170574	3.432	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	453798	15.129	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.469	132	493837	12.527	ng/u1	-0.01
15) 4-Methylphenol-d8	8.675	113	309423	8.007	ng/u1	-0.01
21) Nitrobenzene-d5	9.163	128	321382	17.749	ng/u1	0.00
24) 2-Nitrophenol-d4	9.875	143	353629	17.987	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.405	165	574314	15.462	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.957	131	797413	14.820	ng/u1	-0.01
46) Dimethylphthalate-d6	14.063	166	2068462	17.901	ng/u1	0.00
49) Acenaphthylene-d8	14.345	160	2314101	17.413	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.881	143	63444	3.152	ng/u1	-0.01
60) Fluorene-d10	15.651	176	1830578	18.402	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.792	200	325258	15.869	ng/u1	-0.01
73) Anthracene-d10	17.539	188	3108917	19.276	ng/u1	0.00
81) Pyrene-d10	19.786	212	3698519	19.443	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.974	264	3630596	18.748	ng/u1	-0.01

Target Compounds Qvalue

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

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