

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092923\
 Data File : BP017442.D
 Acq On : 29 Sep 2023 12:48
 Operator : MA/JU
 Sample : 04497-05
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJ38

Quant Time: Sep 29 22:34:30 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	7.957	152	180735	20.000	ng/u1	0.00
20) Naphthalene-d8	10.793	136	754333	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.651	164	486541	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.439	188	1128354	20.000	ng/u1	0.00
79) Chrysene-d12	21.569	240	1098129	20.000	ng/u1	0.00
88) Perylene-d12	24.145	264	1252668	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	10570	2.402	ng/uL	0.00
4) Pyridine-d5	3.740	84	109603	8.910	ng/u1	0.00
7) Phenol-d5	7.110	99	98128	6.611	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.299	67	250667	27.982	ng/u1	0.00
11) 2-Chlorophenol-d4	7.475	132	268954	22.845	ng/u1	0.00
15) 4-Methylphenol-d8	8.675	113	180732	15.661	ng/u1	-0.01
21) Nitrobenzene-d5	9.163	128	175866	32.249	ng/u1	0.00
24) 2-Nitrophenol-d4	9.881	143	187341	31.638	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.410	165	317968	28.423	ng/u1	0.00
31) 4-Chloroaniline-d4	10.963	131	445095	27.466	ng/u1	0.00
46) Dimethylphthalate-d6	14.063	166	1155283	33.146	ng/u1	0.00
49) Acenaphthylene-d8	14.345	160	1315359	32.813	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.892	143	30336	4.996	ng/u1	0.00
60) Fluorene-d10	15.651	176	1033209	34.433	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.792	200	144891	22.783	ng/u1	-0.01
73) Anthracene-d10	17.539	188	1725498	34.480	ng/u1	0.00
81) Pyrene-d10	19.786	212	2102756	35.410	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.974	264	1989977	32.995	ng/u1	-0.01

Target Compounds	Qvalue

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

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