

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
 Data File : BP017330.D
 Acq On : 25 Sep 2023 20:26
 Operator : MA/JU
 Sample : 04450-01
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJ39

Quant Time: Sep 26 00:21:23 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 21 23:51:07 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	203283	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.752	136	895749	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.587	164	572829	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.357	188	1335134	20.000	ng/u1	-0.01
79) Chrysene-d12	21.463	240	1276363	20.000	ng/u1	-0.01
88) Perylene-d12	23.980	264	1423014	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	20201	4.082	ng/uL	0.00
4) Pyridine-d5	3.734	84	201982	14.598	ng/u1	0.00
7) Phenol-d5	7.087	99	120317	7.207	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.269	67	416067	41.294	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.452	132	402036	30.361	ng/u1	-0.01
15) 4-Methylphenol-d8	8.640	113	245182	18.889	ng/u1	-0.02
21) Nitrobenzene-d5	9.128	128	279398	43.145	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.840	143	294015	41.814	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.363	165	488138	36.745	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.916	131	614598	31.938	ng/u1	-0.01
46) Dimethylphthalate-d6	14.004	166	1867012	45.497	ng/u1	-0.01
49) Acenaphthylene-d8	14.287	160	1991671	42.200	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.822	143	40884	5.719	ng/u1	0.00
60) Fluorene-d10	15.587	176	1621141	45.888	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.716	200	269430	35.804	ng/u1	-0.01
73) Anthracene-d10	17.463	188	2808867	47.436	ng/u1	0.00
81) Pyrene-d10	19.704	212	3464875	50.199	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.816	264	3404352	49.689	ng/u1	-0.02

Target Compounds Qvalue

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

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