

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
 Data File : BP017352.D
 Acq On : 26 Sep 2023 09:26
 Operator : MA/JU
 Sample : 04472-12
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EZYK9

Quant Time: Sep 27 00:25:46 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 26 04:52:29 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	190054	20.000	ng/ul	0.00
20) Naphthalene-d8	10.752	136	806063	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.587	164	492547	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.357	188	1109598	20.000	ng/ul	0.00
79) Chrysene-d12	21.463	240	1000598	20.000	ng/ul	0.00
88) Perylene-d12	23.980	264	1111755	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	20981	4.534	ng/uL	0.00
4) Pyridine-d5	3.728	84	295721	22.861	ng/ul	0.00
7) Phenol-d5	7.081	99	392814	25.166	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	249608	26.498	ng/ul	0.00
11) 2-Chlorophenol-d4	7.446	132	312863	25.272	ng/ul	0.00
15) 4-Methylphenol-d8	8.640	113	290752	23.959	ng/ul	0.00
21) Nitrobenzene-d5	9.128	128	160646	27.568	ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143	136711	21.606	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	291296	24.367	ng/ul	0.00
31) 4-Chloroaniline-d4	10.916	131	457926	26.444	ng/ul	0.00
46) Dimethylphthalate-d6	14.004	166	1012793	28.704	ng/ul	0.00
49) Acenaphthylene-d8	14.287	160	1155712	28.479	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143	22273	3.623	ng/ul	0.01
60) Fluorene-d10	15.587	176	903365	29.739	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.716	200	3302	0.528	ng/ul	0.00
73) Anthracene-d10	17.457	188	1396953	28.387	ng/ul	0.00
81) Pyrene-d10	19.704	212	1608819	29.733	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.816	264	1457295	27.225	ng/ul	0.00

Target Compounds	Qvalue

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

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