

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
 Data File : BP017362.D
 Acq On : 26 Sep 2023 15:19
 Operator : MA/JU
 Sample : 04450-06
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBKA2

Quant Time: Sep 27 00:28:46 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	7.922	152	158995	20.000	ng/u1	0.00
20) Naphthalene-d8	10.746	136	693127	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.587	164	421842	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.357	188	935858	20.000	ng/u1	0.00
79) Chrysene-d12	21.463	240	873563	20.000	ng/u1	0.00
88) Perylene-d12	23.974	264	967714	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	17611	4.550	ng/uL	0.00
4) Pyridine-d5	3.734	84	152119	14.057	ng/u1	0.00
7) Phenol-d5	7.081	99	98269	7.525	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	300452	38.126	ng/u1	0.00
11) 2-Chlorophenol-d4	7.446	132	301789	29.139	ng/u1	0.00
15) 4-Methylphenol-d8	8.634	113	192828	18.994	ng/u1	0.00
21) Nitrobenzene-d5	9.122	128	196005	39.116	ng/u1	0.00
24) 2-Nitrophenol-d4	9.834	143	207095	38.063	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	351531	34.197	ng/u1	0.00
31) 4-Chloroaniline-d4	10.910	131	418795	28.125	ng/u1	0.00
46) Dimethylphthalate-d6	13.998	166	1268580	41.979	ng/u1	0.00
49) Acenaphthylene-d8	14.287	160	1366476	39.316	ng/u1	0.00
54) 4-Nitrophenol-d4	14.822	143	25532	4.850	ng/u1	0.00
60) Fluorene-d10	15.581	176	1076069	41.362	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.716	200	138740	26.303	ng/u1	0.00
73) Anthracene-d10	17.457	188	1773137	42.720	ng/u1	0.00
81) Pyrene-d10	19.698	212	2180373	46.155	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.810	264	2140701	45.945	ng/u1	0.00

Target Compounds Qvalue

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

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