

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092923\
 Data File : BP017457.D
 Acq On : 29 Sep 2023 21:51
 Operator : MA/JU
 Sample : 04497-15
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJ60

Quant Time: Sep 29 22:37:44 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	253348	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.793	136	1013689	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.645	164	612311	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.434	188	1289766	20.000	ng/u1	-0.01
79) Chrysene-d12	21.563	240	1270796	20.000	ng/u1	-0.01
88) Perylene-d12	24.133	264	1475351	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	20920	3.392	ng/uL	0.00
4) Pyridine-d5	3.740	84	92239	5.349	ng/u1	0.00
7) Phenol-d5	7.105	99	137207	6.594	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.293	67	321362	25.592	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.469	132	349627	21.186	ng/u1	-0.01
15) 4-Methylphenol-d8	8.669	113	234229	14.479	ng/u1	-0.02
21) Nitrobenzene-d5	9.158	128	208626	28.468	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.875	143	220692	27.735	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.405	165	383192	25.489	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.957	131	448195	20.581	ng/u1	-0.01
46) Dimethylphthalate-d6	14.057	166	1261230	28.753	ng/u1	-0.01
49) Acenaphthylene-d8	14.340	160	1405029	27.850	ng/u1	-0.02
54) 4-Nitrophenol-d4	14.881	143	37870	4.956	ng/u1	-0.01
60) Fluorene-d10	15.645	176	1087597	28.801	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.787	200	161843	22.264	ng/u1	-0.02
73) Anthracene-d10	17.534	188	1767324	30.896	ng/u1	-0.01
81) Pyrene-d10	19.781	212	2096222	30.503	ng/u1	-0.02
92) Benzo(a)pyrene-d12	23.963	264	2147522	30.232	ng/u1	-0.02

Target Compounds	Qvalue

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

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