

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
 Data File : BP017134.D
 Acq On : 13 Sep 2023 01:52
 Operator : MA/JU
 Sample : 04247-16
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJL6

Quant Time: Sep 13 04:11:22 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 12 23:27:27 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	294577	20.000	ng/u1	0.00
20) Naphthalene-d8	10.781	136	1375273	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.610	164	913374	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.381	188	2072705	20.000	ng/u1	0.00
79) Chrysene-d12	21.480	240	1936022	20.000	ng/u1	0.00
88) Perylene-d12	24.004	264	2044893	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.317	96	31823	4.540	ng/uL	0.00
4) Pyridine-d5	3.752	84	192500	9.600	ng/u1	0.00
7) Phenol-d5	7.105	99	157639	6.581	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	334739	22.359	ng/u1	0.00
11) 2-Chlorophenol-d4	7.475	132	215789	11.293	ng/u1	0.00
15) 4-Methylphenol-d8	8.663	113	216646	10.913	ng/u1	0.00
21) Nitrobenzene-d5	9.152	128	202272	21.004	ng/u1	0.00
24) 2-Nitrophenol-d4	9.869	143	101553	9.878	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	209991	10.562	ng/u1	0.00
31) 4-Chloroaniline-d4	10.940	131	523231	17.460	ng/u1	0.00
46) Dimethylphthalate-d6	14.028	166	1144099	17.602	ng/u1	0.00
49) Acenaphthylene-d8	14.310	160	1185798	16.078	ng/u1	0.00
54) 4-Nitrophenol-d4	14.839	143	26579	2.116	ng/u1	0.00
60) Fluorene-d10	15.610	176	888353	15.807	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.734	200	19814	1.711	ng/u1	-0.01
73) Anthracene-d10	17.481	188	1317315	14.959	ng/u1	0.00
81) Pyrene-d10	19.722	212	1539704	15.164	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.839	264	1322680	13.704	ng/u1	-0.01

Target Compounds	Qvalue

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

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