

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092424\
 Data File : BP022006.D
 Acq On : 24 Sep 2024 22:26
 Operator : RC/JU
 Sample : P4092-11
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B47

Quant Time: Sep 24 23:51:55 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 24 15:23:13 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.711	152	100128	20.000	ng/u1	0.00
20) Naphthalene-d8	10.463	136	365281	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.316	164	267767	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.116	188	647373	20.000	ng/u1	0.00
79) Chrysene-d12	21.545	240	739313	20.000	ng/u1	0.00
88) Perylene-d12	24.821	264	897466	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.223	96	11982	4.687	ng/uL	0.00
4) Pyridine-d5	3.640	84	74064	10.136	ng/u1	0.00
7) Phenol-d5	6.893	99	84911	10.367	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	169336	33.575	ng/u1	0.00
11) 2-Chlorophenol-d4	7.252	132	188186	31.756	ng/u1	0.00
15) 4-Methylphenol-d8	8.405	113	147612	22.744	ng/u1	0.00
21) Nitrobenzene-d5	8.846	128	94017	34.388	ng/u1	0.00
24) 2-Nitrophenol-d4	9.558	143	111514	35.973	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.093	165	235474	34.978	ng/u1	0.00
31) 4-Chloroaniline-d4	10.599	131	233121	29.023	ng/u1	0.00
46) Dimethylphthalate-d6	13.722	166	790185	38.176	ng/u1	0.00
49) Acenaphthylene-d8	14.004	160	772993	35.273	ng/u1	0.00
54) 4-Nitrophenol-d4	14.528	143	30032	8.482	ng/u1	0.00
60) Fluorene-d10	15.328	176	664619	36.300	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.440	200	130035	32.354	ng/u1	0.00
73) Anthracene-d10	17.216	188	1181875	39.107	ng/u1	0.00
81) Pyrene-d10	19.598	212	1547810	40.410	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.598	264	1849716	39.186	ng/u1	-0.01

Target Compounds Qvalue

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

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