

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092424\
 Data File : BP022047.D
 Acq On : 26 Sep 2024 04:19
 Operator : RC/JU
 Sample : P4130-06
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0H10

Quant Time: Sep 26 04:49:15 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	7.705	152	88211	20.000	ng/u1	0.00
20) Naphthalene-d8	10.469	136	323336	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.310	164	252637	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.110	188	640594	20.000	ng/u1	0.00
79) Chrysene-d12	21.539	240	782130	20.000	ng/u1	0.00
88) Perylene-d12	24.827	264	926122	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	11233	4.988	ng/uL	0.00
4) Pyridine-d5	3.640	84	76542	11.890	ng/u1	0.00
7) Phenol-d5	6.887	99	70478	9.767	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	126495	28.469	ng/u1	0.00
11) 2-Chlorophenol-d4	7.246	132	146921	28.142	ng/u1	0.00
15) 4-Methylphenol-d8	8.404	113	119521	20.903	ng/u1	0.00
21) Nitrobenzene-d5	8.840	128	75218	31.081	ng/u1	0.00
24) 2-Nitrophenol-d4	9.557	143	88879	32.391	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.093	165	178751	29.997	ng/u1	0.00
31) 4-Chloroaniline-d4	10.598	131	182251	25.633	ng/u1	0.00
46) Dimethylphthalate-d6	13.716	166	644305	32.992	ng/u1	0.00
49) Acenaphthylene-d8	13.998	160	621507	30.059	ng/u1	0.00
54) 4-Nitrophenol-d4	14.522	143	33209	9.941	ng/u1	0.00
60) Fluorene-d10	15.316	176	557540	32.275	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.439	200	127976	32.178	ng/u1	0.00
73) Anthracene-d10	17.204	188	1101950	36.848	ng/u1	-0.01
81) Pyrene-d10	19.586	212	1587464	39.177	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.586	264	1914354	39.300	ng/u1	-0.02

Target Compounds Qvalue

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

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