

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101124\
 Data File : BP022340.D
 Acq On : 11 Oct 2024 18:52
 Operator : RC/JU
 Sample : P4237-12
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4F32

Quant Time: Oct 12 01:06:39 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 08 02:14:47 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.704	152	116211	20.000	ng/u1	0.00
20) Naphthalene-d8	10.469	136	431177	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.322	164	330530	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.110	188	788273	20.000	ng/u1	0.00
79) Chrysene-d12	21.545	240	822017	20.000	ng/u1	0.00
88) Perylene-d12	24.845	264	956681	20.000	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	12194	4.078	ng/uL	0.00
4) Pyridine-d5	3.640	84	163705	19.941	ng/u1	0.00
7) Phenol-d5	6.893	99	228065	23.314	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	134075	23.319	ng/u1	0.00
11) 2-Chlorophenol-d4	7.246	132	178336	25.692	ng/u1	0.00
15) 4-Methylphenol-d8	8.404	113	192155	24.614	ng/u1	0.00
21) Nitrobenzene-d5	8.846	128	86688	26.148	ng/u1	0.00
24) 2-Nitrophenol-d4	9.557	143	105246	27.420	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.098	165	210324	25.782	ng/u1	0.00
31) 4-Chloroaniline-d4	10.592	131	232594	24.129	ng/u1	0.00
46) Dimethylphthalate-d6	13.710	166	717302	27.314	ng/u1	0.00
49) Acenaphthylene-d8	14.010	160	775307	27.796	ng/u1	0.01
54) 4-Nitrophenol-d4	14.545	143	93885	21.310	ng/u1	0.02
60) Fluorene-d10	15.328	176	639211	28.062	ng/u1	0.01
65) 4,6-Dinitro-2-methylph...	15.445	200	99880	19.266	ng/u1	0.00
73) Anthracene-d10	17.216	188	1082093	29.181	ng/u1	0.00
81) Pyrene-d10	19.598	212	1359206	31.268	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.615	264	1526445	29.877	ng/u1	0.02

Target Compounds	Qvalue

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

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