

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
 Data File : BP022382.D
 Acq On : 15 Oct 2024 09:42
 Operator : RC/JU
 Sample : P4349-01
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EOAG9

Quant Time: Oct 15 10:52:36 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Oct 14 11:01:55 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.693	152	113653	20.000	ng/u1	0.00
20) Naphthalene-d8	10.452	136	461908	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.310	164	371914	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.110	188	906967	20.000	ng/u1	0.00
79) Chrysene-d12	21.539	240	963005	20.000	ng/u1	-0.01
88) Perylene-d12	24.845	264	1124560	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.223	96	10666	3.647	ng/uL	0.00
4) Pyridine-d5	3.640	84	51621	6.429	ng/u1	0.00
7) Phenol-d5	6.887	99	63711	6.659	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.034	67	120780	21.480	ng/u1	0.00
11) 2-Chlorophenol-d4	7.240	132	146462	21.575	ng/u1	0.00
15) 4-Methylphenol-d8	8.393	113	116490	15.258	ng/u1	0.00
21) Nitrobenzene-d5	8.840	128	83609	23.541	ng/u1	0.00
24) 2-Nitrophenol-d4	9.552	143	99423	24.180	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.087	165	202590	23.182	ng/u1	0.00
31) 4-Chloroaniline-d4	10.593	131	178430	17.279	ng/u1	0.00
46) Dimethylphthalate-d6	13.710	166	695175	23.526	ng/u1	0.00
49) Acenaphthylene-d8	13.998	160	745572	23.756	ng/u1	0.00
54) 4-Nitrophenol-d4	14.540	143	29635	5.978	ng/u1	0.00
60) Fluorene-d10	15.322	176	637668	24.880	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	132973	22.292	ng/u1	0.00
73) Anthracene-d10	17.222	188	1144444	26.823	ng/u1	0.02
81) Pyrene-d10	19.592	212	1439792	28.272	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.610	264	1643032	27.358	ng/u1	-0.02

Target Compounds Qvalue

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

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