

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
 Data File : BP022404.D
 Acq On : 16 Oct 2024 01:46
 Operator : RC/JU
 Sample : P4349-20
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EOAG2

Quant Time: Oct 16 02:59:43 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	94253	20.000	ng/u1	0.00
20) Naphthalene-d8	10.446	136	348619	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.304	164	271492	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.104	188	660975	20.000	ng/u1	0.00
79) Chrysene-d12	21.539	240	741373	20.000	ng/u1	-0.01
88) Perylene-d12	24.827	264	861284	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	10601	4.371	ng/uL	0.00
4) Pyridine-d5	3.640	84	32202	4.836	ng/u1	0.00
7) Phenol-d5	6.887	99	58866	7.420	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.034	67	134763	28.900	ng/u1	0.00
11) 2-Chlorophenol-d4	7.240	132	154448	27.435	ng/u1	0.00
15) 4-Methylphenol-d8	8.393	113	113393	17.909	ng/u1	0.00
21) Nitrobenzene-d5	8.834	128	86905	32.421	ng/u1	0.00
24) 2-Nitrophenol-d4	9.546	143	103011	33.193	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	208271	31.576	ng/u1	0.00
31) 4-Chloroaniline-d4	10.593	131	55276	7.092	ng/u1	0.00
46) Dimethylphthalate-d6	13.716	166	726643	33.687	ng/u1	0.00
49) Acenaphthylene-d8	13.987	160	751407	32.797	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.534	143	24315	6.719	ng/u1	0.00
60) Fluorene-d10	15.310	176	653096	34.907	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.451	200	135243	31.111	ng/u1	0.00
73) Anthracene-d10	17.204	188	1133624	36.458	ng/u1	0.00
81) Pyrene-d10	19.580	212	1433830	36.572	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.598	264	1711065	37.199	ng/u1	-0.04

Target Compounds Qvalue

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

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