

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100824\
 Data File : BP022272.D
 Acq On : 08 Oct 2024 15:26
 Operator : RC/JU
 Sample : P4162-07
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4ES6

Quant Time: Oct 08 16:20:16 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	7.705	152	97341	20.000	ng/u1	0.00
20) Naphthalene-d8	10.463	136	368845	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.304	164	286960	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.110	188	710946	20.000	ng/u1	0.00
79) Chrysene-d12	21.545	240	907076	20.000	ng/u1	0.00
88) Perylene-d12	24.827	264	1125674	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	7358	2.937	ng/uL	0.00
4) Pyridine-d5	3.640	84	99148	14.418	ng/u1	0.00
7) Phenol-d5	6.893	99	138610	16.916	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	82239	17.077	ng/u1	0.00
11) 2-Chlorophenol-d4	7.252	132	100504	17.286	ng/u1	0.00
15) 4-Methylphenol-d8	8.399	113	117325	17.942	ng/u1	0.00
21) Nitrobenzene-d5	8.846	128	49454	17.438	ng/u1	0.00
24) 2-Nitrophenol-d4	9.557	143	58282	17.750	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.093	165	126208	18.085	ng/u1	0.00
31) 4-Chloroaniline-d4	10.599	131	140858	17.082	ng/u1	0.00
46) Dimethylphthalate-d6	13.716	166	460044	20.178	ng/u1	0.00
49) Acenaphthylene-d8	13.998	160	467926	19.323	ng/u1	0.00
54) 4-Nitrophenol-d4	14.528	143	66931	17.499	ng/u1	0.00
60) Fluorene-d10	15.316	176	402419	20.349	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.457	200	71010	15.187	ng/u1	0.00
73) Anthracene-d10	17.216	188	690353	20.642	ng/u1	0.00
81) Pyrene-d10	19.586	212	925149	19.287	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.604	264	1241613	20.653	ng/u1	0.00

Target Compounds Qvalue

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

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