

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100924\
 Data File : BP022307.D
 Acq On : 10 Oct 2024 03:16
 Operator : RC/JU
 Sample : P4207-14
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EY053

Quant Time: Oct 10 03:52:44 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	111879	20.000	ng/u1	0.00
20) Naphthalene-d8	10.463	136	425660	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.310	164	344949	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.110	188	870957	20.000	ng/u1	0.00
79) Chrysene-d12	21.545	240	908126	20.000	ng/u1	0.00
88) Perylene-d12	24.827	264	1051994	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	9419	3.272	ng/uL	0.00
4) Pyridine-d5	3.646	84	50318	6.366	ng/u1	0.01
7) Phenol-d5	6.893	99	80783	8.578	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	135852	24.543	ng/u1	0.00
11) 2-Chlorophenol-d4	7.246	132	168618	25.233	ng/u1	0.00
15) 4-Methylphenol-d8	8.405	113	142958	19.021	ng/u1	0.00
21) Nitrobenzene-d5	8.846	128	91029	27.813	ng/u1	0.00
24) 2-Nitrophenol-d4	9.557	143	111196	29.346	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.093	165	222443	27.621	ng/u1	0.00
31) 4-Chloroaniline-d4	10.604	131	226872	23.841	ng/u1	0.01
46) Dimethylphthalate-d6	13.716	166	804325	29.347	ng/u1	0.00
49) Acenaphthylene-d8	14.004	160	809464	27.808	ng/u1	0.00
54) 4-Nitrophenol-d4	14.545	143	41489	9.024	ng/u1	0.02
60) Fluorene-d10	15.328	176	719248	30.256	ng/u1	0.01
65) 4,6-Dinitro-2-methylph...	15.445	200	147834	25.808	ng/u1	0.00
73) Anthracene-d10	17.210	188	1432640	34.966	ng/u1	0.00
81) Pyrene-d10	19.586	212	1830414	38.115	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.610	264	1954711	34.793	ng/u1	0.01

Target Compounds Qvalue

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

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