

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121022\
 Data File : BP013046.D
 Acq On : 10 Dec 2022 18:44
 Operator : CG/JU
 Sample : N5950-05
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Dec 12 03:26:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 12 01:32:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	575408	20.000	ng/ul	0.00
20) Naphthalene-d8	10.787	136	2505723	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.610	164	1699238	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.363	188	3922314	20.000	ng/ul	0.00
79) Chrysene-d12	21.451	240	3540051	20.000	ng/ul	0.00
88) Perylene-d12	23.939	264	3422198	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	63586	4.725	ng/uL	0.00
4) Pyridine-d5	3.799	84	37288	0.975	ng/ul	0.02
7) Phenol-d5	7.122	99	317193	6.768	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	886635	31.360	ng/ul	0.00
11) 2-Chlorophenol-d4	7.499	132	849499	23.025	ng/ul	0.00
15) 4-Methylphenol-d8	8.675	113	561947	14.633	ng/ul	0.00
21) Nitrobenzene-d5	9.140	128	612814	33.287	ng/ul	0.00
24) 2-Nitrophenol-d4	9.863	143	647872	31.258	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.399	165	1058948	26.304	ng/ul	0.00
31) 4-Chloroaniline-d4	10.922	131	244744	4.306	ng/ul	0.00
46) Dimethylphthalate-d6	14.016	166	4540638	34.769	ng/ul	0.00
49) Acenaphthylene-d8	14.304	160	4922808	34.305	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	111227	4.735	ng/ul	0.00
60) Fluorene-d10	15.604	176	4016363	36.353	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	722331	28.617	ng/ul	0.00
73) Anthracene-d10	17.463	188	6754062	38.169	ng/ul	0.00
81) Pyrene-d10	19.698	212	7934545	39.048	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.780	264	6634467	38.894	ng/ul	0.00

Target Compounds Qvalue

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

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