

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121022\
 Data File : BP013049.D
 Acq On : 10 Dec 2022 20:26
 Operator : CG/JU
 Sample : N5950-04
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Dec 12 03:27:38 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	595918	20.000	ng/ul	0.00
20) Naphthalene-d8	10.787	136	2634090	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.610	164	1812307	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.363	188	4150511	20.000	ng/ul	0.00
79) Chrysene-d12	21.451	240	3701099	20.000	ng/ul	0.00
88) Perylene-d12	23.939	264	3560899	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	61111	4.385	ng/uL	0.00
4) Pyridine-d5	3.799	84	46114	1.165	ng/ul	0.02
7) Phenol-d5	7.122	99	330181	6.802	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	942219	32.179	ng/ul	0.00
11) 2-Chlorophenol-d4	7.499	132	926542	24.248	ng/ul	0.00
15) 4-Methylphenol-d8	8.675	113	622312	15.647	ng/ul	0.00
21) Nitrobenzene-d5	9.140	128	655098	33.849	ng/ul	0.00
24) 2-Nitrophenol-d4	9.863	143	694878	31.892	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.399	165	1162785	27.476	ng/ul	0.00
31) 4-Chloroaniline-d4	10.922	131	123519	2.067	ng/ul	0.00
46) Dimethylphthalate-d6	14.016	166	4847683	34.804	ng/ul	0.00
49) Acenaphthylene-d8	14.304	160	5047156	32.977	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	135365	5.403	ng/ul	0.00
60) Fluorene-d10	15.604	176	4169102	35.381	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	794802	29.757	ng/ul	0.00
73) Anthracene-d10	17.463	188	7198748	38.445	ng/ul	0.00
81) Pyrene-d10	19.698	212	8433993	39.700	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.780	264	7001334	39.446	ng/ul	0.00

Target Compounds Qvalue

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

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