

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121022\
 Data File : BP013067.D
 Acq On : 11 Dec 2022 07:13
 Operator : CG/JU
 Sample : N5927-10
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Dec 12 03:32:31 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	731500	20.000	ng/ul	0.00
20) Naphthalene-d8	10.781	136	3240990	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.610	164	2192677	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.363	188	5166004	20.000	ng/ul	0.00
79) Chrysene-d12	21.451	240	5054101	20.000	ng/ul	0.00
88) Perylene-d12	23.939	264	5374017	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	70796	4.138	ng/uL	0.00
4) Pyridine-d5	3.781	84	372560	7.666	ng/ul	0.00
7) Phenol-d5	7.122	99	343011	5.757	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	1077660	29.983	ng/ul	0.00
11) 2-Chlorophenol-d4	7.493	132	1052850	22.447	ng/ul	0.00
15) 4-Methylphenol-d8	8.675	113	689933	14.132	ng/ul	0.00
21) Nitrobenzene-d5	9.140	128	716425	30.086	ng/ul	0.00
24) 2-Nitrophenol-d4	9.863	143	771734	28.787	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.398	165	1355373	26.029	ng/ul	0.00
31) 4-Chloroaniline-d4	10.922	131	1740829	23.681	ng/ul	0.00
46) Dimethylphthalate-d6	14.016	166	5759398	34.177	ng/ul	0.00
49) Acenaphthylene-d8	14.304	160	5903574	31.882	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	154878	5.109	ng/ul	0.00
60) Fluorene-d10	15.604	176	5003189	35.094	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.716	200	1030125	30.986	ng/ul	0.00
73) Anthracene-d10	17.463	188	8737441	37.490	ng/ul	0.00
81) Pyrene-d10	19.692	212	10675979	36.800	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.780	264	10406182	38.849	ng/ul	0.00

Target Compounds Qvalue

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

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