

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111022\
 Data File : BP012599.D
 Acq On : 10 Nov 2022 14:57
 Operator : CG/JU
 Sample : N5493-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0AA1

Quant Time: Nov 10 21:35:16 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 09 23:28:28 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	241920	20.000	ng/ul	0.00
20) Naphthalene-d8	10.557	136	1117264	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.416	164	747219	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.181	188	1634630	20.000	ng/ul	0.00
79) Chrysene-d12	21.280	240	1763195	20.000	ng/ul	0.00
88) Perylene-d12	23.651	264	1809933	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	21001	3.538	ng/uL	0.00
4) Pyridine-d5	3.652	84	95226	5.677	ng/ul	0.01
7) Phenol-d5	6.958	99	96883	4.751	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.105	67	341128	27.957	ng/ul	0.00
11) 2-Chlorophenol-d4	7.299	132	310213	19.767	ng/ul	0.00
15) 4-Methylphenol-d8	8.493	113	195104	12.094	ng/ul	0.00
21) Nitrobenzene-d5	8.934	128	222511	26.916	ng/ul	0.00
24) 2-Nitrophenol-d4	9.652	143	222781	24.536	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.199	165	361373	22.092	ng/ul	0.00
31) 4-Chloroaniline-d4	10.716	131	510605	22.160	ng/ul	0.00
46) Dimethylphthalate-d6	13.834	166	1534544	29.795	ng/ul	0.00
49) Acenaphthylene-d8	14.110	160	1652759	28.366	ng/ul	0.00
54) 4-Nitrophenol-d4	14.704	143	26965	3.325	ng/ul	0.04
60) Fluorene-d10	15.416	176	1351444	30.946	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.551	200	228635	25.289	ng/ul	0.00
73) Anthracene-d10	17.281	188	2384727	33.594	ng/ul	0.00
81) Pyrene-d10	19.522	212	2826646	28.608	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.498	264	2944988	33.007	ng/ul	0.00

Target Compounds Qvalue

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

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