

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111022\
 Data File : BP012597.D
 Acq On : 10 Nov 2022 13:48
 Operator : CG/JU
 Sample : N5493-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0AA0

Quant Time: Nov 10 21:34:54 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	264345	20.000	ng/ul	0.00
20) Naphthalene-d8	10.557	136	1215569	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.416	164	813462	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.180	188	1791396	20.000	ng/ul	0.00
79) Chrysene-d12	21.280	240	1948889	20.000	ng/ul	0.00
88) Perylene-d12	23.657	264	2019289	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	22875	3.527	ng/uL	0.00
4) Pyridine-d5	3.652	84	139267	7.599	ng/ul	0.01
7) Phenol-d5	6.957	99	117757	5.284	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.104	67	373952	28.047	ng/ul	0.00
11) 2-Chlorophenol-d4	7.299	132	350926	20.464	ng/ul	0.00
15) 4-Methylphenol-d8	8.493	113	230208	13.059	ng/ul	0.00
21) Nitrobenzene-d5	8.934	128	247871	27.559	ng/ul	0.00
24) 2-Nitrophenol-d4	9.651	143	245620	24.864	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.198	165	407145	22.877	ng/ul	0.00
31) 4-Chloroaniline-d4	10.716	131	560770	22.369	ng/ul	0.00
46) Dimethylphthalate-d6	13.834	166	1597190	28.486	ng/ul	0.00
49) Acenaphthylene-d8	14.110	160	1774577	27.977	ng/ul	0.00
54) 4-Nitrophenol-d4	14.692	143	27836	3.153	ng/ul	0.02
60) Fluorene-d10	15.416	176	1422320	29.917	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.551	200	224591	22.667	ng/ul	0.00
73) Anthracene-d10	17.280	188	2382734	30.629	ng/ul	0.00
81) Pyrene-d10	19.522	212	2862532	26.211	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.504	264	3057559	30.716	ng/ul	0.00

Target Compounds Qvalue

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

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