

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011964.D
 Acq On : 06 Oct 2022 11:37
 Operator : CG/JU
 Sample : N4956-12
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 06 23:04:49 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 05 23:18:55 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	285918	20.000	ng/ul	0.00
20) Naphthalene-d8	10.687	136	1221500	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.522	164	751558	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	1579575	20.000	ng/ul	0.00
79) Chrysene-d12	21.369	240	1518344	20.000	ng/ul	0.00
88) Perylene-d12	23.798	264	1547690	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	26886	4.147	ng/uL	0.00
4) Pyridine-d5	3.728	84	79152	4.168	ng/ul	0.00
7) Phenol-d5	7.040	99	141782	5.744	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	374853	27.398	ng/ul	0.00
11) 2-Chlorophenol-d4	7.411	132	415686	21.333	ng/ul	0.00
15) 4-Methylphenol-d8	8.587	113	266026	13.038	ng/ul	-0.01
21) Nitrobenzene-d5	9.046	128	274897	29.546	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	270565	27.181	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.304	165	475941	25.849	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	508240	18.106	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166	1740474	32.384	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1909838	31.172	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143	48247	4.335	ng/ul	0.00
60) Fluorene-d10	15.522	176	1551580	33.243	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	232259	24.293	ng/ul	0.00
73) Anthracene-d10	17.381	188	2503011	34.787	ng/ul	0.00
81) Pyrene-d10	19.616	212	2918244	37.951	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	2847504	36.754	ng/ul	0.00

Target Compounds Qvalue

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

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