

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011966.D
 Acq On : 06 Oct 2022 12:46
 Operator : CG/JU
 Sample : N4956-08
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 06 23:05:12 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	262968	20.000	ng/ul	0.00
20) Naphthalene-d8	10.687	136	1047167	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.528	164	634882	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	1370905	20.000	ng/ul	0.00
79) Chrysene-d12	21.369	240	1388080	20.000	ng/ul	0.00
88) Perylene-d12	23.804	264	1502093	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	22005	3.691	ng/uL	0.00
4) Pyridine-d5	3.734	84	46169	2.643	ng/ul	0.01
7) Phenol-d5	7.046	99	113340	4.993	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	299589	23.808	ng/ul	0.00
11) 2-Chlorophenol-d4	7.411	132	364038	20.313	ng/ul	0.00
15) 4-Methylphenol-d8	8.593	113	208769	11.124	ng/ul	0.00
21) Nitrobenzene-d5	9.046	128	223735	28.051	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	231640	27.144	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	396896	25.145	ng/ul	0.00
31) 4-Chloroaniline-d4	10.834	131	162908	6.770	ng/ul	0.00
46) Dimethylphthalate-d6	13.940	166	1347379	29.677	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1478917	28.574	ng/ul	0.00
54) 4-Nitrophenol-d4	14.740	143	40853	4.345	ng/ul	0.01
60) Fluorene-d10	15.522	176	1231248	31.228	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	196823	23.720	ng/ul	0.00
73) Anthracene-d10	17.381	188	2030669	32.518	ng/ul	0.00
81) Pyrene-d10	19.616	212	2498988	35.548	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.645	264	2487256	33.078	ng/ul	0.00

Target Compounds Qvalue

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

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