

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011969.D
 Acq On : 06 Oct 2022 14:30
 Operator : CG/JU
 Sample : N4956-14
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 06 23:05:43 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	244933	20.000	ng/ul	0.00
20) Naphthalene-d8	10.687	136	997381	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.528	164	594594	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.286	188	1247009	20.000	ng/ul	0.00
79) Chrysene-d12	21.369	240	1289321	20.000	ng/ul	0.00
88) Perylene-d12	23.798	264	1361457	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	20390	3.671	ng/uL	0.00
4) Pyridine-d5	3.728	84	104590	6.429	ng/ul	0.00
7) Phenol-d5	7.046	99	113747	5.379	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	300709	25.656	ng/ul	0.00
11) 2-Chlorophenol-d4	7.410	132	352034	21.090	ng/ul	0.00
15) 4-Methylphenol-d8	8.593	113	211399	12.094	ng/ul	0.00
21) Nitrobenzene-d5	9.046	128	218168	28.718	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	224468	27.617	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	368022	24.480	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	493784	21.544	ng/ul	0.00
46) Dimethylphthalate-d6	13.939	166	1214721	28.568	ng/ul	0.00
49) Acenaphthylene-d8	14.222	160	1423946	29.376	ng/ul	0.00
54) 4-Nitrophenol-d4	14.745	143	31752	3.606	ng/ul	0.02
60) Fluorene-d10	15.522	176	1133955	30.709	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	172744	22.887	ng/ul	0.00
73) Anthracene-d10	17.380	188	1939741	34.148	ng/ul	0.00
81) Pyrene-d10	19.616	212	2321049	35.546	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.645	264	2385520	35.002	ng/ul	0.00

Target Compounds Qvalue

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

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