

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011980.D
 Acq On : 06 Oct 2022 20:53
 Operator : CG/JU
 Sample : N4923-23
 Misc :
 ALS Vial : 68 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 06 23:08:49 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	7.875	152	286804	20.000	ng/ul	0.00
20) Naphthalene-d8	10.687	136	1191032	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.522	164	708922	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	1485936	20.000	ng/ul	0.00
79) Chrysene-d12	21.369	240	1470291	20.000	ng/ul	0.00
88) Perylene-d12	23.792	264	1489436	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	24401	3.752	ng/uL	0.00
4) Pyridine-d5	3.728	84	84840	4.454	ng/ul	0.00
7) Phenol-d5	7.046	99	93501	3.776	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.205	67	350852	25.564	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.405	132	355647	18.196	ng/ul	0.00
15) 4-Methylphenol-d8	8.587	113	190010	9.283	ng/ul	-0.01
21) Nitrobenzene-d5	9.046	128	239254	26.373	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	230078	23.705	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.304	165	380737	21.208	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	559040	20.426	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166	1334109	26.316	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1535852	26.575	ng/ul	0.00
54) 4-Nitrophenol-d4	14.745	143	22573	2.150	ng/ul	0.02
60) Fluorene-d10	15.516	176	1234205	28.033	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	167419	18.615	ng/ul	0.00
73) Anthracene-d10	17.381	188	2130830	31.480	ng/ul	0.00
81) Pyrene-d10	19.616	212	2455893	32.982	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	2390101	32.056	ng/ul	0.00

Target Compounds Qvalue

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

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