

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011931.D
 Acq On : 05 Oct 2022 15:39
 Operator : CG/JU
 Sample : N4890-09
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EY0E5

Quant Time: Oct 05 23:22:55 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	254564	20.000	ng/u1	0.00
20) Naphthalene-d8	10.693	136	1037165	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.528	164	609745	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.286	188	1276561	20.000	ng/u1	0.00
79) Chrysene-d12	21.368	240	1309540	20.000	ng/u1	0.00
88) Perylene-d12	23.798	264	1327966	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	24503	4.245	ng/uL	0.00
4) Pyridine-d5	3.728	84	76212	4.508	ng/u1	0.00
7) Phenol-d5	7.046	99	126653	5.763	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	368173	30.224	ng/u1	0.00
11) 2-Chlorophenol-d4	7.410	132	405963	23.401	ng/u1	0.00
15) 4-Methylphenol-d8	8.593	113	217068	11.949	ng/u1	0.00
21) Nitrobenzene-d5	9.051	128	248409	31.444	ng/u1	0.00
24) 2-Nitrophenol-d4	9.769	143	253376	29.978	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	423585	27.095	ng/u1	0.00
31) 4-Chloroaniline-d4	10.834	131	359006	15.063	ng/u1	0.00
46) Dimethylphthalate-d6	13.939	166	1446202	33.167	ng/u1	0.00
49) Acenaphthylene-d8	14.222	160	1659086	33.377	ng/u1	0.00
54) 4-Nitrophenol-d4	14.739	143	34023	3.768	ng/u1	0.01
60) Fluorene-d10	15.522	176	1329022	35.097	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	188367	24.379	ng/u1	0.00
73) Anthracene-d10	17.386	188	1939242	33.349	ng/u1	0.00
81) Pyrene-d10	19.621	212	2451095	36.958	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.645	264	2436025	36.645	ng/u1	0.00

Target Compounds	Qvalue

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

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