

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
 Data File : BP011930.D
 Acq On : 05 Oct 2022 15:04
 Operator : CG/JU
 Sample : N4890-12
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EY0E7

Quant Time: Oct 05 23:22:44 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	240354	20.000	ng/u1	0.00
20) Naphthalene-d8	10.693	136	974399	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.528	164	581882	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.286	188	1243225	20.000	ng/u1	0.00
79) Chrysene-d12	21.369	240	1278808	20.000	ng/u1	0.00
88) Perylene-d12	23.804	264	1290533	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	30086	5.521	ng/uL	0.00
4) Pyridine-d5	3.728	84	128094	8.024	ng/u1	0.00
7) Phenol-d5	7.046	99	137817	6.642	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	395484	34.386	ng/u1	0.00
11) 2-Chlorophenol-d4	7.410	132	441158	26.933	ng/u1	0.00
15) 4-Methylphenol-d8	8.593	113	258046	15.044	ng/u1	0.00
21) Nitrobenzene-d5	9.052	128	277718	37.419	ng/u1	0.00
24) 2-Nitrophenol-d4	9.775	143	273022	34.383	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	453528	30.879	ng/u1	0.00
31) 4-Chloroaniline-d4	10.834	131	632279	28.238	ng/u1	0.00
46) Dimethylphthalate-d6	13.939	166	1580202	37.976	ng/u1	0.00
49) Acenaphthylene-d8	14.222	160	1812692	38.213	ng/u1	0.00
54) 4-Nitrophenol-d4	14.739	143	35884	4.164	ng/u1	0.01
60) Fluorene-d10	15.522	176	1456327	40.301	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	210376	27.957	ng/u1	0.00
73) Anthracene-d10	17.386	188	2395069	42.292	ng/u1	0.00
81) Pyrene-d10	19.622	212	2884968	44.546	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.645	264	2834616	43.878	ng/u1	0.00

Target Compounds Qvalue

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

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