

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011898.D
 Acq On : 04 Oct 2022 06:26
 Operator : CG/JU
 Sample : N4859-16
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E10048

Quant Time: Oct 04 06:56:53 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 01 00:53:59 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	242175	20.000	ng/u1	0.00
20) Naphthalene-d8	10.699	136	973923	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.534	164	561932	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.286	188	1171722	20.000	ng/u1	-0.01
79) Chrysene-d12	21.375	240	1272390	20.000	ng/u1	-0.01
88) Perylene-d12	23.810	264	1326549	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	24609	4.482	ng/uL	0.00
4) Pyridine-d5	3.734	84	47353	2.944	ng/u1	0.00
7) Phenol-d5	7.052	99	158270	7.570	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	364298	31.436	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.416	132	418322	25.347	ng/u1	0.00
15) 4-Methylphenol-d8	8.599	113	290040	16.782	ng/u1	0.00
21) Nitrobenzene-d5	9.052	128	259904	35.036	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.775	143	262551	33.081	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.316	165	432409	29.455	ng/u1	0.00
31) 4-Chloroaniline-d4	10.840	131	490517	21.917	ng/u1	0.00
46) Dimethylphthalate-d6	13.945	166	1473103	36.659	ng/u1	0.00
49) Acenaphthylene-d8	14.228	160	1660038	36.238	ng/u1	0.00
54) 4-Nitrophenol-d4	14.740	143	46275	5.561	ng/u1	0.00
60) Fluorene-d10	15.528	176	1335546	38.270	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	209906	29.597	ng/u1	0.00
73) Anthracene-d10	17.386	188	2183945	40.918	ng/u1	-0.01
81) Pyrene-d10	19.622	212	2742777	42.564	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.651	264	2907090	43.778	ng/u1	-0.01

Target Compounds Qvalue

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

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