

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082322\
 Data File : BP011543.D
 Acq On : 23 Aug 2022 15:53
 Operator : CG/JU
 Sample : SP5970
 Misc : SFAM SURROGATE
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SP5970

Quant Time: Aug 23 16:39:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 17 01:58:37 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.552	152	66422	20.000	ng/u1	0.00
20) Naphthalene-d8	10.340	136	281728	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.228	164	161592	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.998	188	308184	20.000	ng/u1	0.00
79) Chrysene-d12	21.133	240	276173	20.000	ng/u1	0.00
88) Perylene-d12	23.427	264	279508	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.064	96	34182	15.970	ng/uL	0.00
4) Pyridine-d5	3.469	84	430416	78.542	ng/u1	-0.01
7) Phenol-d5	6.740	99	446724	73.187	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.905	67	298355	75.236	ng/u1	0.00
11) 2-Chlorophenol-d4	7.087	132	373260	83.143	ng/u1	0.00
15) 4-Methylphenol-d8	8.275	113	367809	75.285	ng/u1	0.00
21) Nitrobenzene-d5	8.722	128	182623	90.040	ng/u1	0.00
24) 2-Nitrophenol-d4	9.440	143	194372	94.164	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.969	165	312065	81.901	ng/u1	0.00
31) 4-Chloroaniline-d4	10.498	131	495861	78.178	ng/u1	0.00
46) Dimethylphthalate-d6	13.657	166	969357	84.814	ng/u1	0.00
49) Acenaphthylene-d8	13.922	160	1132801	89.178	ng/u1	0.00
54) 4-Nitrophenol-d4	14.463	143	160599	72.932	ng/u1	0.00
60) Fluorene-d10	15.234	176	789357	82.094	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.375	200	135613	81.436	ng/u1	0.00
73) Anthracene-d10	17.104	188	1191402	87.288	ng/u1	0.00
81) Pyrene-d10	19.369	212	1267810	90.734	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.280	264	1177428	86.599	ng/u1	0.00

Target Compounds Qvalue

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

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