

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072622\
 Data File : BP011106.D
 Acq On : 26 Jul 2022 12:45
 Operator : CG/JU
 Sample : N3801-17
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBHY2

Quant Time: Jul 26 14:44:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 26 14:31:36 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.664	152	505427	20.000	ng/u1	0.00
20) Naphthalene-d8	10.458	136	2319778	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.328	164	1364351	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.092	188	2441843	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.210	240	1478385	20.000	ng/u1	# 0.00
88) Perylene-d12	23.551	264	1439496	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	59426	4.239	ng/uL	0.00
4) Pyridine-d5	3.576	84	285417	7.876	ng/u1	0.00
7) Phenol-d5	6.840	99	290869	7.011	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.005	67	924496	34.653	ng/u1	0.00
11) 2-Chlorophenol-d4	7.193	132	869798	25.945	ng/u1	0.00
15) 4-Methylphenol-d8	8.375	113	595573	17.931	ng/u1	0.00
21) Nitrobenzene-d5	8.828	128	580760	34.354	ng/u1	0.00
24) 2-Nitrophenol-d4	9.546	143	580709	34.715	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	944880	30.887	ng/u1	0.00
31) 4-Chloroaniline-d4	10.605	131	919651	18.235	ng/u1	0.00
46) Dimethylphthalate-d6	13.751	166	3474669	37.262	ng/u1	0.00
49) Acenaphthylene-d8	14.016	160	3999434	36.351	ng/u1	0.00
54) 4-Nitrophenol-d4	14.546	143	62534	3.314	ng/u1	0.00
60) Fluorene-d10	15.328	176	2748243	34.673	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	396511	32.816	ng/u1	0.00
73) Anthracene-d10	17.192	188	3765034	36.269	ng/u1	0.00
81) Pyrene-d10	19.451	212	3535208	45.530	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.404	264	2588589	37.441	ng/u1	0.00

Target Compounds Qvalue

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

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