

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072522\
 Data File : BP011095.D
 Acq On : 25 Jul 2022 16:45
 Operator : CG/JU
 Sample : N3801-06
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBHX1

Quant Time: Jul 25 23:21:34 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 19 03:25:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.664	152	530103	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.458	136	2474609	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.328	164	1574541	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.092	188	3158031	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.210	240	2296662	20.000	ng/u1	#-0.02
88) Perylene-d12	23.551	264	1868976	20.000	ng/u1	-0.03

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	57760	3.928	ng/uL	-0.02
4) Pyridine-d5	3.576	84	366162	9.634	ng/u1	-0.02
7) Phenol-d5	6.846	99	299428	6.881	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.011	67	1057529	37.794	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.199	132	974379	27.712	ng/u1	-0.01
15) 4-Methylphenol-d8	8.381	113	628911	18.054	ng/u1	-0.02
21) Nitrobenzene-d5	8.828	128	720240	39.939	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.546	143	707988	39.676	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.081	165	1087127	33.313	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.605	131	1739464	32.333	ng/u1	-0.01
46) Dimethylphthalate-d6	13.751	166	4536609	42.156	ng/u1	-0.01
49) Acenaphthylene-d8	14.022	160	5287101	41.640	ng/u1	0.00
54) 4-Nitrophenol-d4	14.546	143	82352	3.781	ng/u1	-0.01
60) Fluorene-d10	15.328	176	3726122	40.735	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.457	200	575468	36.826	ng/u1	0.00
73) Anthracene-d10	17.198	188	5497069	40.945	ng/u1	0.00
81) Pyrene-d10	19.451	212	5773179	47.862	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.404	264	3743737	41.706	ng/u1	-0.02

Target Compounds	Qvalue

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

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