

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072122\
 Data File : BP011068.D
 Acq On : 21 Jul 2022 18:29
 Operator : CG/JU
 Sample : N3709-13
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Jul 21 23:25:09 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	7.675	152	410215	20.000	ng/ul	0.00
20) Naphthalene-d8	10.463	136	1726379	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.333	164	993696	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.104	188	1958344	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.227	240	1792645	20.000	ng/ul	# 0.00
88) Perylene-d12	23.580	264	1834315	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	12066	1.060	ng/uL	0.00
4) Pyridine-d5	3.575	84	678	0.023	ng/ul	-0.02
7) Phenol-d5	6.852	99	366922	10.897	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.016	67	262019	12.101	ng/ul	0.00
11) 2-Chlorophenol-d4	7.204	132	326136	11.986	ng/ul	0.00
15) 4-Methylphenol-d8	8.387	113	237355	8.805	ng/ul	-0.01
21) Nitrobenzene-d5	8.834	128	156828	12.465	ng/ul	0.00
24) 2-Nitrophenol-d4	9.551	143	133817	10.749	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.087	165	257496	11.310	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.610	131	171581	4.572	ng/ul	0.00
46) Dimethylphthalate-d6	13.751	166	624646	9.197	ng/ul	-0.01
49) Acenaphthylene-d8	14.022	160	1132507	14.133	ng/ul	0.00
54) 4-Nitrophenol-d4	14.557	143	40405	2.940	ng/ul	0.00
60) Fluorene-d10	15.333	176	833348	14.436	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.457	200	24211	2.498	ng/ul	0.00
73) Anthracene-d10	17.198	188	1304231	15.666	ng/ul	0.00
81) Pyrene-d10	19.457	212	1619214	17.198	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.421	264	1499195	17.017	ng/ul	0.00

Target Compounds	Qvalue

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

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