

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072122\
 Data File : BP011061.D
 Acq On : 21 Jul 2022 14:26
 Operator : CG/JU
 Sample : N3709-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Jul 21 23:23:49 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	403075	20.000	ng/u1	0.00
20) Naphthalene-d8	10.463	136	1652759	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.333	164	926715	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.104	188	1861426	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.227	240	1762498	20.000	ng/u1	# 0.00
88) Perylene-d12	23.580	264	1814607	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	70214	6.280	ng/uL	0.00
4) Pyridine-d5	3.587	84	934904	32.351	ng/u1	0.00
7) Phenol-d5	6.851	99	1116318	33.739	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.016	67	752093	35.349	ng/u1	0.00
11) 2-Chlorophenol-d4	7.204	132	921965	34.485	ng/u1	0.00
15) 4-Methylphenol-d8	8.387	113	897048	33.866	ng/u1	-0.01
21) Nitrobenzene-d5	8.834	128	443963	36.860	ng/u1	0.00
24) 2-Nitrophenol-d4	9.551	143	419135	35.168	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.087	165	740019	33.953	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.610	131	1175578	32.717	ng/u1	0.00
46) Dimethylphthalate-d6	13.757	166	2353109	37.152	ng/u1	0.00
49) Acenaphthylene-d8	14.028	160	2933557	39.255	ng/u1	0.00
54) 4-Nitrophenol-d4	14.551	143	290520	22.664	ng/u1	0.00
60) Fluorene-d10	15.333	176	2013081	37.392	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.457	200	179719	19.512	ng/u1	0.00
73) Anthracene-d10	17.204	188	3030346	38.294	ng/u1	0.00
81) Pyrene-d10	19.462	212	3449210	37.262	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.427	264	3300336	37.868	ng/u1	0.00

Target Compounds	Qvalue

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

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