

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072622\
 Data File : BP011111.D
 Acq On : 26 Jul 2022 15:38
 Operator : CG/JU
 Sample : N3801-11
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBHX4

Quant Time: Jul 26 16:55:05 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	7.663	152	416200	20.000	ng/u1	0.00
20) Naphthalene-d8	10.457	136	1860537	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.328	164	1026325	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.092	188	1733721	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.210	240	1239072	20.000	ng/u1	# 0.00
88) Perylene-d12	23.557	264	1323704	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	53804	4.660	ng/uL	0.00
4) Pyridine-d5	3.575	84	300262	10.062	ng/u1	0.00
7) Phenol-d5	6.840	99	234151	6.854	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.005	67	792702	36.083	ng/u1	0.00
11) 2-Chlorophenol-d4	7.193	132	727836	26.365	ng/u1	0.00
15) 4-Methylphenol-d8	8.381	113	475787	17.396	ng/u1	0.00
21) Nitrobenzene-d5	8.828	128	495148	36.519	ng/u1	0.00
24) 2-Nitrophenol-d4	9.546	143	481478	35.887	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	761319	31.029	ng/u1	0.00
31) 4-Chloroaniline-d4	10.604	131	1135229	28.066	ng/u1	0.00
46) Dimethylphthalate-d6	13.751	166	2751644	39.228	ng/u1	0.00
49) Acenaphthylene-d8	14.016	160	3216111	38.859	ng/u1	0.00
54) 4-Nitrophenol-d4	14.545	143	37551	2.645	ng/u1	0.00
60) Fluorene-d10	15.328	176	2190300	36.735	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	276566	32.238	ng/u1	0.00
73) Anthracene-d10	17.192	188	2908788	39.465	ng/u1	0.00
81) Pyrene-d10	19.451	212	2862745	43.991	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.404	264	2570908	40.438	ng/u1	0.00

Target Compounds Qvalue

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

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