

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072522\
 Data File : BP011094.D
 Acq On : 25 Jul 2022 16:10
 Operator : CG/JU
 Sample : N3801-16
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBHY1

Quant Time: Jul 25 23:21:25 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	7.663	152	489633	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.457	136	2269362	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.328	164	1435592	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.092	188	2805779	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.210	240	1856302	20.000	ng/u1	#-0.02
88) Perylene-d12	23.551	264	1563492	20.000	ng/u1	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	55696	4.101	ng/uL	-0.02
4) Pyridine-d5	3.581	84	182532	5.200	ng/u1	-0.01
7) Phenol-d5	6.846	99	252889	6.292	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.011	67	867302	33.558	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.199	132	823734	25.364	ng/u1	-0.01
15) 4-Methylphenol-d8	8.381	113	499071	15.510	ng/u1	-0.02
21) Nitrobenzene-d5	8.828	128	574355	34.729	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.546	143	581866	35.557	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.081	165	890809	29.766	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.604	131	969996	19.661	ng/u1	-0.01
46) Dimethylphthalate-d6	13.751	166	3749581	38.215	ng/u1	-0.01
49) Acenaphthylene-d8	14.022	160	4235285	36.585	ng/u1	0.00
54) 4-Nitrophenol-d4	14.545	143	70638	3.557	ng/u1	-0.01
60) Fluorene-d10	15.328	176	3043521	36.493	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.457	200	475593	34.255	ng/u1	0.00
73) Anthracene-d10	17.192	188	4221460	35.391	ng/u1	-0.01
81) Pyrene-d10	19.451	212	4455887	45.705	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.404	264	2860334	38.090	ng/u1	-0.02

Target Compounds Qvalue

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

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