

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072821\
 Data File : BP006401.D
 Acq On : 28 Jul 2021 21:29
 Operator : CG/JU
 Sample : M3107-11
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGD01

Quant Time: Jul 29 01:08:14 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 24 04:29:12 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	314709	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.563	136	1383475	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.410	164	795327	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.157	188	1553873	20.000	ng/u1	-0.01
79) Chrysene-d12	21.257	240	1476260	20.000	ng/u1	-0.01
88) Perylene-d12	23.598	264	1458340	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	46204	5.480	ng/uL	0.00
4) Pyridine-d5	3.699	84	648281	25.901	ng/u1	0.00
7) Phenol-d5	6.958	99	871713	29.493	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	568926	30.935	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.316	132	663499	30.215	ng/u1	-0.01
15) 4-Methylphenol-d8	8.487	113	717340	30.862	ng/u1	-0.01
21) Nitrobenzene-d5	8.934	128	336839	32.699	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.652	143	341023	34.671	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.187	165	585720	30.388	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.704	131	945416	29.794	ng/u1	-0.01
46) Dimethylphthalate-d6	13.828	166	1961124	34.646	ng/u1	-0.01
49) Acenaphthylene-d8	14.098	160	2364044	33.993	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.610	143	330612	29.416	ng/u1	0.00
60) Fluorene-d10	15.404	176	1613202	33.901	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.522	200	218397	25.212	ng/u1	-0.01
73) Anthracene-d10	17.257	188	2464560	35.136	ng/u1	-0.01
81) Pyrene-d10	19.504	212	2754255	36.273	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.451	264	2572645	34.714	ng/u1	-0.01

Target Compounds Qvalue

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

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