

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031122\
 Data File : BN018911.D
 Acq On : 10 Mar 2022 20:13
 Operator : CG/JU
 Sample : N1814-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DTW-3

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 03/11/2022
 Supervised By : Jagrut Upadhyay 03/11/2022

Quant Time: Mar 11 03:05:20 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN030322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 11 03:02:57 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	4949	0.400	ng	0.00
7) Naphthalene-d8	10.712	136	14965	0.400	ng	# 0.00
13) Acenaphthene-d10	14.538	164	10223	0.400	ng	0.00
19) Phenanthrene-d10	17.277	188	21034	0.400	ng	0.00
29) Chrysene-d12	21.449	240	20589	0.400	ng	# 0.00
35) Perylene-d12	23.840	264	19044	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.464	112	1784	0.165	ng	0.00
5) Phenol-d6	7.060	99	1430	0.110	ng	0.00
8) Nitrobenzene-d5	9.057	82	3055m	0.280	ng	0.00
11) 2-Methylnaphthalene-d10	12.299	152	7512	0.308	ng	0.00
14) 2,4,6-Tribromophenol	16.015	330	2100	0.446	ng	-0.01
15) 2-Fluorobiphenyl	13.169	172	13126	0.300	ng	0.00
27) Fluoranthene-d10	19.303	212	24264	0.392	ng	0.00
31) Terphenyl-d14	19.893	244	20459	0.465	ng	0.00
Target Compounds						Qvalue
24) Pentachlorophenol	16.928	266	290	0.263	ng	92
25) Phenanthrene	17.318	178	1642	0.024	ng	# 83
34) Bis(2-ethylhexyl)phtha...	21.360	149	6035	0.126	ng	# 98

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

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