

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090421\
 Data File : BN016229.D
 Acq On : 05 Sep 2021 15:49
 Operator : CG/JU
 Sample : M3561-06
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 VE4-11

Quant Time: Sep 06 06:23:48 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN090421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 06 02:08:58 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.873	152	14128	0.400	ng	0.00
7) Naphthalene-d8	10.660	136	72606	0.400	ng	# 0.00
13) Acenaphthene-d10	14.497	164	57216	0.400	ng	0.00
19) Phenanthrene-d10	17.249	188	100757	0.400	ng	# 0.00
29) Chrysene-d12	21.450	240	104882	0.400	ng	# 0.02
35) Perylene-d12	23.854	264	77944	0.400	ng	# 0.03
System Monitoring Compounds						
4) 2-Fluorophenol	5.457	112	5192	0.130	ng	0.00
5) Phenol-d6	7.043	99	3853	0.076	ng	0.00
8) Nitrobenzene-d5	9.025	82	21422	0.368	ng	0.00
11) 2-Methylnaphthalene-d10	12.252	152	40318	0.349	ng	0.00
14) 2,4,6-Tribromophenol	15.994	330	7989	0.313	ng	0.00
15) 2-Fluorobiphenyl	13.114	172	57015	0.264	ng	-0.01
27) Fluoranthene-d10	19.296	212	87463	0.289	ng	0.01
31) Terphenyl-d14	19.892	244	78623	0.297	ng	0.02
Target Compounds						
						Qvalue
9) Naphthalene	10.711	128	6633	0.035	ng	# 35
12) 2-Methylnaphthalene	12.323	142	6320	0.049	ng	# 24
16) Acenaphthylene	14.218	152	21685	0.084	ng	# 1
17) Acenaphthene	14.561	154	3809	0.023	ng	# 51
25) Phenanthrene	17.285	178	18003	0.064	ng	# 1
26) Anthracene	17.383	178	125294	0.466	ng	# 84
28) Fluoranthene	19.326	202	22212	0.066	ng	# 85
30) Pyrene	19.693	202	61990	0.181	ng	# 86
32) Benzo(a)anthracene	21.440	228	15543	0.044	ng	# 1
33) Chrysene	21.493	228	25326	0.074	ng	# 1
34) Bis(2-ethylhexyl)phtha...	21.355	149	42526	0.179	ng	# 92
36) Indeno(1,2,3-cd)pyrene	26.336	276	7574	0.035	ng	# 91
37) Benzo(b)fluoranthene	23.122	252	13873	0.052	ng	# 1
38) Benzo(k)fluoranthene	23.170	252	13734	0.056	ng	# 1
39) Benzo(a)pyrene	23.748	252	11337	0.048	ng	# 1
40) Dibenzo(a,h)anthracene	26.349	278	6248	0.035	ng	# 1
41) Benzo(g,h,i)perylene	27.096	276	5684	0.031	ng	# 1

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

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