

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031921\
 Data File : BN013985.D
 Acq On : 18 Mar 2021 16:55
 Operator : CG/JU
 Sample : M1714-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GWBR3-209-219-031721

Quant Time: Mar 18 17:29:37 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN022421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 18 15:58:40 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.708	152	5992	0.40	ng	0.00
7) Naphthalene-d8	10.492	136	22952	0.40	ng	# 0.00
13) Acenaphthene-d10	14.346	164	12927	0.40	ng	0.00
19) Phenanthrene-d10	17.092	188	29135	0.40	ng	# 0.00
29) Chrysene-d12	21.271	240	29531	0.40	ng	# 0.00
36) Perylene-d12	23.499	264	26542	0.40	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.300	112	2221	0.14	ng	0.00
5) Phenol-d6	6.871	99	1683	0.08	ng	0.00
8) Nitrobenzene-d5	8.857	82	4602	0.31	ng	0.00
11) 2-Methylnaphthalene-d10	12.080	152	10357	0.28	ng	0.00
14) 2,4,6-Tribromophenol	15.831	330	1460	0.28	ng	0.00
15) 2-Fluorobiphenyl	12.963	172	15737	0.34	ng	0.00
27) Fluoranthene-d10	19.119	212	31083	0.36	ng	0.00
31) Terphenyl-d14	19.720	244	41220	0.63	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.247	88	31990	4.42	ng	96
9) Naphthalene	10.530	128	1585	0.03	ng	# 85
12) 2-Methylnaphthalene	12.155	142	4947	0.12	ng	98
25) Phenanthrene	17.128	178	3762	0.05	ng	98
28) Fluoranthene	19.149	202	4706	0.05	ng	# 88
30) Pyrene	19.511	202	3902	0.04	ng	# 95
32) Benzo(a)anthracene	21.250	228	2298	0.03	ng	# 90
33) Chrysene	21.303	228	3080	0.03	ng	# 94
34) Bis(2-ethylhexyl)phtha...	21.186	149	3840	0.09	ng	98
37) Benzo(b)fluoranthene	22.832	252	3817	0.04	ng	# 47
38) Benzo(k)fluoranthene	22.877	252	2072	0.03	ng	# 36
39) Benzo(a)pyrene	23.400	252	2110	0.03	ng	# 12
41) Benzo(g,h,i)perylene	26.419	276	1740	0.02	ng	# 80

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

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