

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031921\
 Data File : BN013987.D
 Acq On : 18 Mar 2021 18:07
 Operator : CG/JU
 Sample : M1714-03DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

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

Quant Time: Mar 18 18:37:11 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	6369	0.40	ng	0.00
7) Naphthalene-d8	10.492	136	23898	0.40	ng	# 0.00
13) Acenaphthene-d10	14.346	164	13054	0.40	ng	0.00
19) Phenanthrene-d10	17.092	188	28530	0.40	ng	# 0.00
29) Chrysene-d12	21.271	240	27767	0.40	ng	# 0.00
36) Perylene-d12	23.499	264	25021	0.40	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.300	112	426	0.02	ng	0.00
5) Phenol-d6	6.871	99	333	0.01	ng	0.00
8) Nitrobenzene-d5	8.857	82	822	0.05	ng	0.00
11) 2-Methylnaphthalene-d10	12.084	152	1871	0.05	ng	0.00
14) 2,4,6-Tribromophenol	15.831	330	236	0.04	ng	0.00
15) 2-Fluorobiphenyl	12.963	172	2862	0.06	ng	0.00
27) Fluoranthene-d10	19.119	212	5139	0.06	ng	0.00
31) Terphenyl-d14	19.719	244	6914	0.11	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.247	88	6065	0.79	ng	94
12) 2-Methylnaphthalene	12.155	142	1051	0.03	ng	# 94
28) Fluoranthene	19.149	202	2321	0.03	ng	# 95
30) Pyrene	19.511	202	2257	0.03	ng	98
32) Benzo(a)anthracene	21.250	228	2205	0.03	ng	94
33) Chrysene	21.303	228	2464	0.03	ng	94
34) Bis(2-ethylhexyl)phtha...	21.186	149	1377	0.03	ng	# 97
37) Benzo(b)fluoranthene	22.832	252	2601	0.03	ng	# 52
38) Benzo(k)fluoranthene	22.876	252	2060	0.03	ng	# 48
39) Benzo(a)pyrene	23.404	252	1744	0.02	ng	# 32
41) Benzo(g,h,i)perylene	26.419	276	1621	0.02	ng	# 82

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

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