

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070221\
 Data File : BN015314.D
 Acq On : 03 Jul 2021 01:57
 Operator : CG/JU
 Sample : M2898-03
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW-10B-062921

Quant Time: Jul 03 03:03:27 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN063021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 01 09:06:31 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.734	152	21830	0.40	ng	# 0.00
7) Naphthalene-d8	10.495	136	99462	0.40	ng	0.00
13) Acenaphthene-d10	14.345	164	53045	0.40	ng	0.00
19) Phenanthrene-d10	17.090	188	104135	0.40	ng	# 0.00
29) Chrysene-d12	21.291	240	78591	0.40	ng	-0.01
35) Perylene-d12	23.532	264	63059	0.40	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.356	112	5384	0.10	ng	0.00
5) Phenol-d6	6.905	99	3933	0.06	ng	0.00
8) Nitrobenzene-d5	8.860	82	17992	0.32	ng	0.00
11) 2-Methylnaphthalene-d10	12.083	152	44581	0.27	ng	0.00
14) 2,4,6-Tribromophenol	15.842	330	4841	0.43	ng	0.00
15) 2-Fluorobiphenyl	12.962	172	70861	0.33	ng	0.00
27) Fluoranthene-d10	19.127	212	117350	0.38	ng	0.00
31) Terphenyl-d14	19.728	244	102835	0.56	ng	-0.01
Target Compounds						
2) 1,4-Dioxane	3.345	88	52276	5.11	ng	# 61
24) Pentachlorophenol	16.750	266	53	0.25	ng	# 86
32) Benzo(a)anthracene	21.270	228	12959	0.05	ng	98
33) Chrysene	21.323	228	13232	0.05	ng	98
34) Bis(2-ethylhexyl)phtha...	21.206	149	55014	0.70	ng	# 87
36) Indeno(1,2,3-cd)pyrene	25.791	276	6698	0.03	ng	# 87
37) Benzo(b)fluoranthene	22.855	252	11525	0.05	ng	# 53
38) Benzo(k)fluoranthene	22.902	252	10849	0.04	ng	# 85
39) Benzo(a)pyrene	23.433	252	8800	0.04	ng	# 1
40) Dibenzo(a,h)anthracene	25.805	278	5133	0.03	ng	# 79
41) Benzo(g,h,i)perylene	26.476	276	6230	0.03	ng	# 82

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

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