

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100421\
 Data File : BN016762.D
 Acq On : 06 Oct 2021 03:53
 Operator : CG/JU
 Sample : M3915-18
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RW4-20210922

Quant Time: Oct 06 04:32:22 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 04 17:52:33 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.642	152	25771	0.400	ng	0.00
7) Naphthalene-d8	10.407	136	114398	0.400	ng	# 0.00
13) Acenaphthene-d10	14.269	164	59896	0.400	ng	0.00
19) Phenanthrene-d10	17.030	188	114900	0.400	ng	0.00
29) Chrysene-d12	21.227	240	115743	0.400	ng	#-0.01
35) Perylene-d12	23.484	264	117433	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.273	112	8681	0.132	ng	0.02
5) Phenol-d6	6.831	99	5104	0.070	ng	0.00
8) Nitrobenzene-d5	8.784	82	19771	0.246	ng	0.00
11) 2-Methylnaphthalene-d10	12.003	152	55296	0.324	ng	0.00
14) 2,4,6-Tribromophenol	15.778	330	6875	0.316	ng	0.00
15) 2-Fluorobiphenyl	12.886	172	70040	0.292	ng	-0.01
27) Fluoranthene-d10	19.068	212	128720	0.409	ng	0.00
31) Terphenyl-d14	19.678	244	102959	0.398	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.253	88	51696	4.298	ng	# 69
24) Pentachlorophenol	16.689	266	150	0.342	ng	92
34) Bis(2-ethylhexyl)phtha...	21.164	149	21978	0.101	ng	99

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

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