

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101321\
 Data File : BN016943.D
 Acq On : 14 Oct 2021 00:42
 Operator : CG/JU
 Sample : M4063-01RE
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 P287-DITCHRE

Manual Integrations
 APPROVED

mohammad
 10/18/2021 9:03:18 AM

Quant Time: Oct 14 02:11:44 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN101321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 13 02:19:57 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.688	152	21289	0.400	ng	0.00
7) Naphthalene-d8	10.445	136	97294	0.400	ng	#-0.01
13) Acenaphthene-d10	14.294	164	53553	0.400	ng	0.00
19) Phenanthrene-d10	17.030	188	103724	0.400	ng	#-0.01
29) Chrysene-d12	21.227	240	104179	0.400	ng	0.00
35) Perylene-d12	23.471	264	101602	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.310	112	7559	0.156	ng	0.02
5) Phenol-d6	6.877	99	5289	0.098	ng	0.00
8) Nitrobenzene-d5	8.822	82	18947	0.306	ng	0.00
11) 2-Methylnaphthalene-d10	12.038	152	50404	0.370	ng	0.00
14) 2,4,6-Tribromophenol	15.791	330	11033	0.454	ng	0.00
15) 2-Fluorobiphenyl	12.924	172	69484	0.343	ng	0.00
27) Fluoranthene-d10	19.068	212	115307	0.443	ng	0.00
31) Terphenyl-d14	19.678	244	102551	0.448	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.290	88	374m	0.026	ng	Qvalue
24) Pentachlorophenol	16.689	266	676	0.217	ng	97
34) Bis(2-ethylhexyl)phtha...	21.164	149	21434	0.299	ng	99

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

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