

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101321\
 Data File : BN016916.D
 Acq On : 13 Oct 2021 02:37
 Operator : CG/JU
 Sample : M4170-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 P287-DITCH

Manual Integrations
 APPROVED

mohammad
 10/14/2021 10:52:06 AM

Quant Time: Oct 13 06:19:49 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	7.689	152	20912	0.40	ng	0.00
7) Naphthalene-d8	10.445	136	94082	0.40	ng	#-0.01
13) Acenaphthene-d10	14.294	164	52413	0.40	ng	0.00
19) Phenanthrene-d10	17.030	188	99178	0.40	ng	#-0.01
29) Chrysene-d12	21.228	240	98638	0.40	ng	# 0.00
35) Perylene-d12	23.467	264	102952	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.310	112	7409	0.16	ng	0.02
5) Phenol-d6	6.877	99	4861	0.09	ng	0.00
8) Nitrobenzene-d5	8.823	82	18850	0.32	ng	0.00
11) 2-Methylnaphthalene-d10	12.038	152	44994	0.34	ng	0.00
14) 2,4,6-Tribromophenol	15.791	330	10407	0.44	ng	0.00
15) 2-Fluorobiphenyl	12.924	172	61799	0.31	ng	0.00
27) Fluoranthene-d10	19.068	212	107256	0.43	ng	0.00
31) Terphenyl-d14	19.679	244	89413	0.41	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.290	88	526m	0.04	ng	
24) Pentachlorophenol	16.689	266	200	0.20	ng	99
34) Bis(2-ethylhexyl)phtha...	21.164	149	17391	0.28	ng	99

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

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