

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101321\
 Data File : BN016918.D
 Acq On : 13 Oct 2021 03:48
 Operator : CG/JU
 Sample : M4154-05
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SW-CB-32

Manual Integrations
 APPROVED

mohammad
 10/14/2021 10:52:08 AM

Quant Time: Oct 13 06:20:03 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	21346	0.40	ng	0.00
7) Naphthalene-d8	10.445	136	93332	0.40	ng	#-0.01
13) Acenaphthene-d10	14.294	164	55052	0.40	ng	0.00
19) Phenanthrene-d10	17.030	188	105706	0.40	ng	#-0.01
29) Chrysene-d12	21.227	240	105669	0.40	ng	# 0.00
35) Perylene-d12	23.474	264	114087	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.309	112	7440	0.15	ng	0.02
5) Phenol-d6	6.877	99	4710	0.09	ng	0.00
8) Nitrobenzene-d5	8.822	82	19285	0.32	ng	0.00
11) 2-Methylnaphthalene-d10	12.038	152	37578	0.29	ng	0.00
14) 2,4,6-Tribromophenol	15.791	330	10031	0.40	ng	0.00
15) 2-Fluorobiphenyl	12.924	172	50177	0.24	ng	0.00
27) Fluoranthene-d10	19.068	212	87973	0.33	ng	0.00
31) Terphenyl-d14	19.678	244	69143	0.30	ng	0.00
Target Compounds						Qvalue
24) Pentachlorophenol	16.689	266	4168	0.32	ng	100
25) Phenanthrene	17.078	178	29188m	0.10	ng	

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

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