

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121720\
 Data File : BM028143.D
 Acq On : 17 Dec 2020 16:20
 Operator : JU/CG
 Sample : L5088-13
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TR3-A-D-COMP

Quant Time: Dec 17 16:51:48 2020
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM121620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 16 01:58:38 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.210	152	113226	20.00	ng	0.00
21) Naphthalene-d8	11.040	136	423389	20.00	ng	# 0.00
39) Acenaphthene-d10	14.834	164	216043	20.00	ng	-0.01
64) Phenanthrene-d10	17.557	188	417191	20.00	ng	0.00
76) Chrysene-d12	21.668	240	415350	20.00	ng	# 0.00
86) Perylene-d12	24.039	264	458908	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.711	112	831419	121.01	ng	0.00
7) Phenol-d6	7.352	99	895571	91.28	ng	0.00
23) Nitrobenzene-d5	9.387	82	751228	88.05	ng	-0.01
42) 2,4,6-Tribromophenol	16.310	330	347308	121.10	ng	0.00
45) 2-Fluorobiphenyl	13.469	172	1475955	88.82	ng	0.00
79) Terphenyl-d14	20.127	244	1749141	76.41	ng	0.00

Target Compounds	Qvalue

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

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