

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080921\
 Data File : BF125019.D
 Acq On : 10 Aug 2021 00:50
 Operator : JU/CG
 Sample : M3308-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 72-11930

Manual Integrations
 APPROVED

mohammad
 8/10/2021 4:35:05 PM

Quant Time: Aug 10 04:47:29 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 09 11:03:44 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.822	152	5181	20.000	ng	0.00
21) Naphthalene-d8	8.104	136	19889	20.000	ng	# 0.00
39) Acenaphthene-d10	9.945	164	16898m	20.000	ng	0.08
64) Phenanthrene-d10	11.527	188	6579m	20.000	ng	0.18
76) Chrysene-d12	14.062	240	8843m	20.000	ng	0.08
86) Perylene-d12	15.474	264	6953	20.000	ng	0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	30669	96.968	ng	0.00
7) Phenol-d6	6.463	99	38030	95.360	ng	0.00
23) Nitrobenzene-d5	7.386	82	23184	60.979	ng	0.00
42) 2,4,6-Tribromophenol	10.804	330	5576m	36.935	ng	0.15
45) 2-Fluorobiphenyl	9.192	172	47900m	41.478	ng	0.01
79) Terphenyl-d14	13.092	244	23511	44.848	ng	0.16
Target Compounds						
84) Bis(2-ethylhexyl)phtha...	14.009	149	6568m	15.274	ng	Qvalue

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

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