

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082520\
 Data File : BF121425.D
 Acq On : 25 Aug 2020 13:52
 Operator : JU/CG
 Sample : L3784-16
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RO-04

Manual Integrations
 APPROVED

mohammad
 8/26/2020 4:00:28 PM

Quant Time: Aug 25 14:35:51 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF081920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 20 04:56:10 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	151920	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	597367	20.00	ng	0.00
39) Acenaphthene-d10	9.73	164	313314	20.00	ng	0.00
64) Phenanthrene-d10	11.21	188	623103	20.00	ng	0.00
76) Chrysene-d12	13.85	240	574733	20.00	ng	0.00
86) Perylene-d12	15.27	264	595087	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	685330	73.97	ng	0.01
7) Phenol-d6	6.35	99	845541	71.49	ng	0.00
23) Nitrobenzene-d5	7.26	82	551500	52.21	ng	0.00
42) 2,4,6-Tribromophenol	10.52	330	258745	70.87	ng	0.00
45) 2-Fluorobiphenyl	9.05	172	958152	53.99	ng	0.00
79) Terphenyl-d14	12.80	244	1018211	34.62	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.45	163	297954	13.071	ng	99
71) Phenanthrene	11.23	178	77899	2.416	ng	99
75) Fluoranthene	12.42	202	149609	4.274	ng	98
78) Pyrene	12.65	202	129419	2.998	ng	99
88) Benzo(b)fluoranthene	14.86	252	92651m	2.410	ng	

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

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