

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071619\
 Data File : BF115607.D
 Acq On : 16 Jul 2019 19:44
 Operator : HP/JU
 Sample : K3826-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RBR-200034

Manual Integrations
 APPROVED

mohammad
 7/17/2019 12:06:35 PM

Quant Time: Jul 17 05:30:33 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 14:03:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	171723	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	648702	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	333693	20.00	ng	-0.01
64) Phenanthrene-d10	11.40	188	604312	20.00	ng	0.00
76) Chrysene-d12	14.04	240	402411	20.00	ng	-0.01
87) Perylene-d12	15.51	264	470323	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	1272226	121.18	ng	0.00
7) Phenol-d6	6.52	99	1653349	124.11	ng	0.00
23) Nitrobenzene-d5	7.44	82	1055951	94.65	ng	-0.01
42) 2,4,6-Tribromophenol	10.70	330	433217	111.22	ng	-0.01
45) 2-Fluorobiphenyl	9.24	172	1662221	87.19	ng	-0.01
79) Terphenyl-d14	12.99	244	1616850	74.14	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.63	163	360321	14.352	ng	99
71) Phenanthrene	11.43	178	130581	4.215	ng	95
75) Fluoranthene	12.61	202	168056	5.134	ng	98
78) Pyrene	12.84	202	132705	3.892	ng	96
81) Benzo(a)anthracene	14.03	228	58170	2.194	ng	99
83) Chrysene	14.06	228	66595	2.502	ng	97
88) Benzo(b)fluoranthene	15.07	252	76022m	2.510	ng	

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

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