

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071619\
 Data File : BF115605.D
 Acq On : 16 Jul 2019 18:51
 Operator : HP/JU
 Sample : K3830-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RT-3456

Manual Integrations
 APPROVED

mohammad
 7/17/2019 12:06:33 PM

Quant Time: Jul 17 05:21:36 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	152217	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	562095	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	275755	20.00	ng	-0.01
64) Phenanthrene-d10	11.40	188	504736	20.00	ng	0.00
76) Chrysene-d12	14.04	240	356468	20.00	ng	-0.01
87) Perylene-d12	15.51	264	429974	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1003475	107.83	ng	0.00
7) Phenol-d6	6.51	99	1230175	104.18	ng	0.00
23) Nitrobenzene-d5	7.44	82	676199	69.95	ng	-0.01
42) 2,4,6-Tribromophenol	10.70	330	276640	85.95	ng	-0.01
45) 2-Fluorobiphenyl	9.24	172	1296566	82.30	ng	-0.01
79) Terphenyl-d14	12.99	244	1223996	63.36	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.63	163	208648	10.057	ng	98
71) Phenanthrene	11.42	178	56544	2.186	ng	94
75) Fluoranthene	12.61	202	100010	3.658	ng	98
78) Pyrene	12.84	202	104441	3.458	ng	97
83) Chrysene	14.06	228	58741	2.492	ng	96
88) Benzo(b)fluoranthene	15.07	252	70649m	2.552	ng	
90) Benzo(a)pyrene	15.44	252	47872	2.012	ng	99

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

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