

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071919\
 Data File : BF115673.D
 Acq On : 19 Jul 2019 13:27
 Operator : HP/JU
 Sample : K3874-01 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HJDC-GRABA

Manual Integrations
 APPROVED

mohammad
 7/22/2019 10:01:27 AM

Quant Time: Jul 19 15:24:47 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	184291	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	744319	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	347060	20.00	ng	-0.01
64) Phenanthrene-d10	11.40	188	599751	20.00	ng	0.00
76) Chrysene-d12	14.04	240	356431	20.00	ng	0.00
87) Perylene-d12	15.52	264	405916	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	446861	39.66	ng	0.00
7) Phenol-d6	6.50	99	600204	41.98	ng	-0.01
23) Nitrobenzene-d5	7.43	82	368410	28.78	ng	-0.02
42) 2,4,6-Tribromophenol	10.70	330	140325	34.64	ng	-0.01
45) 2-Fluorobiphenyl	9.23	172	623667	31.45	ng	-0.02
79) Terphenyl-d14	12.99	244	482289	24.97	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.63	163	99732	3.819	ng	98
71) Phenanthrene	11.43	178	466980	15.190	ng	96
72) Anthracene	11.47	178	114390	3.600	ng	97
75) Fluoranthene	12.62	202	401707	12.365	ng	99
78) Pyrene	12.84	202	318241	10.538	ng	96
81) Benzo(a)anthracene	14.03	228	139151	5.926	ng	97
83) Chrysene	14.07	228	122647	5.203	ng	97
86) Indeno(1,2,3-cd)pyrene	16.99	276	54556	2.724	ng	94
88) Benzo(b)fluoranthene	15.09	252	145864m	5.581	ng	
90) Benzo(a)pyrene	15.45	252	95578	4.255	ng	99
92) Benzo(g,h,i)perylene	17.44	276	49629	2.523	ng	99

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

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