

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070519\
 Data File : BF115395.D
 Acq On : 5 Jul 2019 18:09
 Operator : HP/JU
 Sample : K3622-03 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-02-070319-B

Manual Integrations
 APPROVED

mohammad
 7/8/2019 2:13:03 PM

Quant Time: Jul 06 05:50:44 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 05 14:57:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	164135	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	657958	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	321858	20.00	ng	0.00
64) Phenanthrene-d10	11.41	188	558398	20.00	ng	0.00
76) Chrysene-d12	14.05	240	409611	20.00	ng	-0.02
87) Perylene-d12	15.53	264	392458	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	500906	47.15	ng	0.00
7) Phenol-d6	6.51	99	677674	50.16	ng	-0.01
23) Nitrobenzene-d5	7.45	82	395903	35.55	ng	-0.01
42) 2,4,6-Tribromophenol	10.72	330	168288	47.29	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	792356	43.20	ng	0.00
79) Terphenyl-d14	12.99	244	750420	37.51	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.64	163	90524	3.747	ng	99
71) Phenanthrene	11.43	178	240843	8.201	ng	97
72) Anthracene	11.49	178	72495	2.462	ng	98
75) Fluoranthene	12.62	202	342793	11.086	ng	97
78) Pyrene	12.85	202	335033	10.518	ng	98
81) Benzo(a)anthracene	14.04	228	173344	6.605	ng	94
83) Chrysene	14.07	228	156340	5.798	ng	95
88) Benzo(b)fluoranthene	15.09	252	194747m	7.621	ng	
89) Benzo(k)fluoranthene	15.12	252	73547m	3.236	ng	
90) Benzo(a)pyrene	15.46	252	134331	6.095	ng	97
92) Benzo(g,h,i)perylene	17.44	276	64210	3.602	ng	96

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

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