

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF042219\
 Data File : BF113898.D
 Acq On : 23 Apr 2019 00:28
 Operator : JU/SJ
 Sample : K2202-03 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SA-06-041819-B

Manual Integrations
 APPROVED

Sohil
 4/23/2019 4:08:14 PM

Quant Time: Apr 23 06:03:33 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF042219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 23 03:32:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	176188	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	601537	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	295272	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	627279	20.00	ng	0.00
76) Chrysene-d12	14.06	240	592315	20.00	ng	-0.01
87) Perylene-d12	15.54	264	477292	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	523148	50.78	ng	0.00
7) Phenol-d6	6.52	99	644019	49.83	ng	0.00
23) Nitrobenzene-d5	7.45	82	359510	36.90	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	196323	54.58	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	640346	33.92	ng	0.00
79) Terphenyl-d14	13.00	244	867660	30.03	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.64	163	72808	3.409	ng	Qvalue 100
71) Phenanthrene	11.44	178	96667	3.066	ng	98
75) Fluoranthene	12.63	202	187483	5.451	ng	98
78) Pyrene	12.86	202	207144	5.001	ng	98
81) Benzo(a)anthracene	14.04	228	116902	3.350	ng	96
83) Chrysene	14.08	228	98299	2.892	ng	98
88) Benzo(b)fluoranthene	15.10	252	110583m	3.662	ng	
90) Benzo(a)pyrene	15.47	252	84357	3.230	ng	# 91

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

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