

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF041919\
 Data File : BF113861.D
 Acq On : 19 Apr 2019 16:29
 Operator : JU/SJ
 Sample : K2202-01 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 4/22/2019 7:26:46 PM

Quant Time: Apr 22 03:16:20 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF041819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 18 13:55:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	205523	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	731844	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	336788	20.00	ng	0.00
64) Phenanthrene-d10	11.43	188	631462	20.00	ng	0.02
76) Chrysene-d12	14.43	240	550278m	20.00	ng	0.39
87) Perylene-d12	16.32	264	484014m	20.00	ng	0.81
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	636005	51.46	ng	0.01
7) Phenol-d6	6.52	99	809766	52.53	ng	0.00
23) Nitrobenzene-d5	7.46	82	448266	37.30	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	213691	57.43	ng	0.01
45) 2-Fluorobiphenyl	9.26	172	819495	37.11	ng	0.00
79) Terphenyl-d14	13.14	244	890617	33.94	ng	0.14
Target Compounds						
50) Dimethylphthalate	9.65	163	138961	5.896	ng	100
71) Phenanthrene	11.45	178	132910	4.133	ng	97
75) Fluoranthene	12.72	202	303566	8.906	ng	97
78) Pyrene	12.97	202	294178	7.586	ng	97
81) Benzo(a)anthracene	14.42	228	146936m	4.568	ng	
83) Chrysene	14.46	228	141153m	4.421	ng	
86) Indeno(1,2,3-cd)pyrene	17.88	276	68376m	2.623	ng	
88) Benzo(b)fluoranthene	15.82	252	172086m	5.703	ng	
90) Benzo(a)pyrene	16.24	252	116394m	4.363	ng	
92) Benzo(g,h,i)perylene	18.34	276	63039m	2.613	ng	

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

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