

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110118\
 Data File : BF110378.D
 Acq On : 1 Nov 2018 20:07
 Operator : JU/SJ
 Sample : J5192-05 2X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-05-103118-A

Manual Integrations
 APPROVED

Sohil
 11/2/2018 2:48:47 PM

Quant Time: Nov 02 04:53:47 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 16:06:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	75497	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	281197	20.00	ng	0.00
39) Acenaphthene-d10	10.00	164	106740	20.00	ng	0.00
64) Phenanthrene-d10	11.49	188	172623	20.00	ng	0.00
76) Chrysene-d12	14.13	240	182917	20.00	ng	0.00
87) Perylene-d12	15.66	264	184749	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	207040	44.66	ng	0.00
7) Phenol-d6	6.57	99	248197	41.85	ng	0.00
23) Nitrobenzene-d5	7.52	82	151146	31.88	ng	0.00
42) 2,4,6-Tribromophenol	10.79	330	39565	37.42	ng	0.00
45) 2-Fluorobiphenyl	9.31	172	258687	38.06	ng	0.00
79) Terphenyl-d14	13.07	244	218355	24.83	ng	0.00
Target Compounds						
10) Phenol	6.59	94	14569	2.298	ng	Qvalue # 67
50) Dimethylphthalate	9.70	163	31537	4.003	ng	97
71) Phenanthrene	11.51	178	24713	2.736	ng	98
75) Fluoranthene	12.70	202	54905	5.989	ng	99
78) Pyrene	12.93	202	52374	3.994	ng	97
81) Benzo(a)anthracene	14.12	228	31798	2.931	ng	95
83) Chrysene	14.16	228	26674	2.589	ng	97
88) Benzo(b)fluoranthene	15.20	252	36196m	3.393	ng	
90) Benzo(a)pyrene	15.59	252	25775	2.626	ng	99

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

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