

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110218\
 Data File : BF110432.D
 Acq On : 3 Nov 2018 00:37
 Operator : JU/SJ
 Sample : J5769-11
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FLOOD-AREA-2

Manual Integrations
 APPROVED

Sohil
 11/5/2018 10:32:53 AM

Quant Time: Nov 03 03:22:59 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	77428	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	284214	20.00	ng	0.00
39) Acenaphthene-d10	10.00	164	109249	20.00	ng	0.00
64) Phenanthrene-d10	11.49	188	188939	20.00	ng	0.00
76) Chrysene-d12	14.14	240	133711	20.00	ng	0.00
87) Perylene-d12	15.66	264	99347	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	476028	100.12	ng	0.00
7) Phenol-d6	6.58	99	588748	96.81	ng	0.00
23) Nitrobenzene-d5	7.52	82	342405	71.46	ng	0.00
42) 2,4,6-Tribromophenol	10.79	330	101367	93.67	ng	0.00
45) 2-Fluorobiphenyl	9.31	172	602293	86.57	ng	0.00
79) Terphenyl-d14	13.07	244	492105	76.54	ng	0.00
Target Compounds						
10) Phenol	6.59	94	38311	5.893	ng	71
50) Dimethylphthalate	9.70	163	91590	11.357	ng	99
71) Phenanthrene	11.51	178	120036	12.142	ng	97
72) Anthracene	11.57	178	34353	3.467	ng	97
75) Fluoranthene	12.71	202	444463	44.299	ng	98
78) Pyrene	12.94	202	396724	41.386	ng	99
81) Benzo(a)anthracene	14.13	228	215633	27.194	ng	98
83) Chrysene	14.16	228	232296	30.844	ng	98
86) Indeno(1,2,3-cd)pyrene	17.23	276	89254	14.285	ng	99
88) Benzo(b)fluoranthene	15.21	252	272461m	47.492	ng	
89) Benzo(k)fluoranthene	15.23	252	79801m	14.149	ng	
90) Benzo(a)pyrene	15.60	252	154275	29.225	ng	99
91) Dibenzo(a,h)anthracene	17.23	278	25054m	5.500	ng	
92) Benzo(a,h,i)perylene	17.71	276	83844	18.680	ng	99

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

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