

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050118\
 Data File : BF104992.D
 Acq On : 1 May 2018 19:01
 Operator : JU/SJ
 Sample : J2590-03
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EPE-MX-4(0-5)

Manual Integrations
 APPROVED

Sohil
 5/2/2018 4:21:34 PM

Quant Time: May 02 11:47:22 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 01 16:13:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	118441	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	495100	20.00	ng	0.00
38) Acenaphthene-d10	9.83	164	222943	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	368915	20.00	ng	0.00
75) Chrysene-d12	14.02	240	230914	20.00	ng	0.04
86) Perylene-d12	15.56	264	213443	20.00	ng	0.13
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	648643	83.74	ng	0.00
7) Phenol-d6	6.43	99	811155	85.23	ng	0.00
23) Nitrobenzene-d5	7.36	82	499710	65.91	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	195328	84.46	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	910881	64.54	ng	0.00
78) Terphenyl-d14	12.92	244	751871	74.23	ng	0.00
Target Compounds						
10) Phenol	6.44	94	39649	3.926	ng	81
49) Dimethylphthalate	9.55	163	134736	7.664	ng	100
74) Fluoranthene	12.54	202	72823	3.393	ng	97
77) Pyrene	12.77	202	72223	4.220	ng	98
80) Benzo(a)anthracene	14.00	228	34722	2.517	ng	94
82) Chrysene	14.04	228	33224m	2.345	ng	
87) Benzo(b)fluoranthene	15.13	252	46586m	3.456	ng	
91) Benzo(g,h,i)perylene	17.49	276	19416	2.023	ng	97

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

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