

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070617\
 Data File : BF096539.D
 Acq On : 6 Jul 2017 17:59
 Operator : SJ/MA
 Sample : I3999-02
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS2-1-4-COMP

Manual Integrations
 APPROVED

mohammad
 7/7/2017 1:30:12 PM

Quant Time: Jul 07 03:23:01 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 05 13:27:10 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	114584	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	486522	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	188705	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	267563	20.00	ng	0.00
75) Chrysene-d12	13.86	240	249895	20.00	ng	0.00
86) Perylene-d12	15.28	264	188801	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	832415	107.22	ng	0.01
7) Phenol-d6	6.36	99	1060209	111.09	ng	0.00
23) Nitrobenzene-d5	7.27	82	566119	69.92	ng	0.00
41) 2,4,6-Tribromophenol	10.54	330	146593	99.45	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	983872	89.14	ng	0.00
78) Terphenyl-d14	12.82	244	617932	52.84	ng	0.00
Target Compounds						
10) Phenol	6.37	94	27269	2.508	ng	Qvalue 90
49) Dimethylphthalate	9.47	163	281603	19.990	ng	98
70) Phenanthrene	11.25	178	33767	2.335	ng	96
74) Fluoranthene	12.44	202	93720	6.609	ng	99
77) Pyrene	12.66	202	89110	4.443	ng	98
80) Benzo(a)anthracene	13.85	228	49299	3.407	ng	97
82) Chrysene	13.89	228	46203	3.049	ng	97
87) Benzo(b)fluoranthene	14.87	252	50455m	4.265	ng	
89) Benzo(a)pyrene	15.22	252	33104	3.134	ng	# 76

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

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