

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112316\
 Data File : BF091292.D
 Acq On : 23 Nov 2016 23:15
 Operator : UM/SJ
 Sample : H5740-08
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-03(0-0.16)

Manual Integrations
 APPROVED

sohil
 11/28/2016 5:18:14 PM

Quant Time: Nov 24 00:48:50 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 23 13:22:53 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	220807	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	920718	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	456261	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	696170	20.00	ng	0.00
75) Chrysene-d12	14.03	240	508950	20.00	ng	-0.01
86) Perylene-d12	15.47	264	491008	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	1858140	135.21	ng	0.02
7) Phenol-d6	6.51	99	2152246	121.89	ng	0.01
23) Nitrobenzene-d5	7.45	82	1565441	97.88	ng	0.00
41) 2,4,6-Tribromophenol	10.70	330	538470	117.03	ng	0.00
44) 2-Fluorobiphenyl	9.24	172	2195985	78.81	ng	0.00
78) Terphenyl-d14	12.99	244	1441165	64.67	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.52	94	127815	6.49	ng	83
32) Benzoic acid	7.89	122	95222	8.49	ng	96
49) Dimethylphthalate	9.63	163	323931	11.37	ng	96
74) Fluoranthene	12.61	202	117237	3.19	ng	95
77) Pyrene	12.83	202	109263	2.91	ng	97
80) Benzo(a)anthracene	14.02	228	80307	2.78	ng	97
82) Chrysene	14.05	228	58793m	2.15	ng	
85) Indeno(1,2,3-cd)pyrene	16.88	276	48823	2.08	ng	94
87) Benzo(b)fluoranthene	15.06	252	117075m	3.78	ng	
89) Benzo(a)pyrene	15.41	252	71939m	2.73	ng	

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

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