

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061417\
 Data File : BF095944.D
 Acq On : 15 Jun 2017 2:45
 Operator : SJ/MA
 Sample : I3665-04
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HA3-WC1

Manual Integrations
 APPROVED

mohammad
 6/15/2017 3:55:55 PM

Quant Time: Jun 15 04:18:19 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 05 16:15:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.33	152	71572	20.00	ng	-0.06
21) Naphthalene-d8	9.37	136	278143	20.00	ng	-0.06
38) Acenaphthene-d10	12.19	164	109984	20.00	ng	-0.07
63) Phenanthrene-d10	14.58	188	164908	20.00	ng	-0.06
75) Chrysene-d12	18.30	240	137372	20.00	ng	-0.05
86) Perylene-d12	19.97	264	99576	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	455465	101.40	ng	-0.06
7) Phenol-d6	6.90	99	545805	101.50	ng	-0.05
23) Nitrobenzene-d5	8.25	82	305059	68.11	ng	-0.07
41) 2,4,6-Tribromophenol	13.49	330	84428	86.76	ng	-0.06
44) 2-Fluorobiphenyl	11.15	172	523671	66.26	ng	-0.07
78) Terphenyl-d14	16.96	244	324216	58.57	ng	-0.06
Target Compounds						
10) Phenol	6.93	94	21428	3.84	ng	Qvalue # 64
49) Dimethylphthalate	11.84	163	70651	8.46	ng	98
51) Acenaphthene	12.24	154	15882	2.27	ng	94
57) Fluorene	13.08	166	22973	2.68	ng	# 92
70) Phenanthrene	14.62	178	116476	12.24	ng	97
71) Anthracene	14.71	178	43177	4.35	ng	97
74) Fluoranthene	16.40	202	174064	18.48	ng	98
77) Pyrene	16.70	202	154904	15.11	ng	95
80) Benzo(a)anthracene	18.29	228	65741	8.23	ng	97
82) Chrysene	18.34	228	57160	7.16	ng	99
85) Indeno(1,2,3-cd)pyrene	21.15	276	18434	2.70	ng	92
87) Benzo(b)fluoranthene	19.57	252	57260m	9.18	ng	
88) Benzo(k)fluoranthene	19.60	252	18188m	3.05	ng	
89) Benzo(a)pyrene	19.92	252	40543	7.07	ng	# 94
91) Benzo(a,h,i)perylene	21.49	276	16510	3.33	ng	98

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

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