

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062516\
 Data File : BF088377.D
 Acq On : 25 Jun 2016 15:18
 Operator : UM/SJ
 Sample : H3622-10
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SC-STA-1456(0-4)

Quant Time: Jun 26 02:46:52 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 22 14:55:14 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.56	152	79505	20.00	ng	-0.05
21) Naphthalene-d8	7.85	136	307217	20.00	ng	-0.05
38) Acenaphthene-d10	9.59	164	135517	20.00	ng	-0.06
63) Phenanthrene-d10	11.06	188	228075	20.00	ng	-0.06
75) Chrysene-d12	13.69	240	158599	20.00	ng	-0.06
86) Perylene-d12	15.05	264	143100	20.00	ng	-0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.14	112	456646	98.19	ng	-0.03
7) Phenol-d6	6.23	99	562880	99.87	ng	-0.03
23) Nitrobenzene-d5	7.14	82	357772	68.00	ng	-0.03
41) 2,4,6-Tribromophenol	10.39	330	118657	82.92	ng	-0.05
44) 2-Fluorobiphenyl	8.92	172	656738	75.51	ng	-0.05
78) Terphenyl-d14	12.65	244	469683	67.39	ng	-0.05
Target Compounds						Qvalue
49) Dimethylphthalate	9.32	163	112762	11.49	ng	99
74) Fluoranthene	12.27	202	99926	7.91	ng	96
77) Pyrene	12.49	202	102101	8.68	ng	97
80) Benzo(a)anthracene	13.68	228	62624	6.93	ng	96
82) Chrysene	13.71	228	48890m	5.75	ng	
85) Indeno(1,2,3-cd)pyrene	16.29	276	25420	3.22	ng	# 100
87) Benzo(b)fluoranthene	14.69	252	77090m	7.73	ng	
88) Benzo(k)fluoranthene	14.70	252	38993m	5.18	ng	
89) Benzo(a)pyrene	14.99	252	44163	5.47	ng	95
91) Benzo(g,h,i)perylene	16.66	276	20604	2.67	ng	97

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

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