

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042816\
 Data File : BF086484.D
 Acq On : 29 Apr 2016 1:15
 Operator : UM/SJ
 Sample : H2654-13
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 26WARD-2A-CS-19(0-6FTBGS)

Manual Integrations
 APPROVED

sohil
 4/29/2016 4:52:54 PM

Quant Time: Apr 29 06:55:31 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 27 14:20:52 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.78	152	143374	20.00	ng	-0.02
21) Naphthalene-d8	8.08	136	602503	20.00	ng	-0.02
38) Acenaphthene-d10	9.82	164	242622	20.00	ng	-0.02
63) Phenanthrene-d10	11.31	188	357892	20.00	ng	-0.01
75) Chrysene-d12	13.94	240	304834	20.00	ng	-0.02
86) Perylene-d12	15.35	264	281246	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	843739	101.51	ng	-0.01
7) Phenol-d6	6.43	99	1243948	107.23	ng	-0.01
23) Nitrobenzene-d5	7.36	82	831075	74.35	ng	-0.02
41) 2,4,6-Tribromophenol	10.61	330	148306	57.96	ng	-0.02
44) 2-Fluorobiphenyl	9.15	172	1133810	80.93	ng	-0.02
78) Terphenyl-d14	12.89	244	760969	55.00	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.55	163	171374	9.38	ng	99
70) Phenanthrene	11.33	178	98587	5.25	ng	98
74) Fluoranthene	12.51	202	143814	7.59	ng	98
77) Pyrene	12.74	202	134788	6.14	ng	98
80) Benzo(a)anthracene	13.93	228	80116	4.93	ng	98
82) Chrysene	13.96	228	60886m	3.45	ng	
85) Indeno(1,2,3-cd)pyrene	16.68	276	41029	3.13	ng	96
87) Benzo(b)fluoranthene	14.95	252	90382m	5.46	ng	
88) Benzo(k)fluoranthene	14.96	252	46975m	2.72	ng	
89) Benzo(a)pyrene	15.29	252	67585	4.33	ng	# 92
91) Benzo(g,h,i)perylene	17.09	276	37705	2.94	ng	96

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

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