

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF093015\
 Data File : BF081926.D
 Acq On : 1 Oct 2015 3:25
 Operator : UM/IZ
 Sample : G3828-02
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OGS-SSCF-PH-FL-2-3.5-4

Manual Integrations
 APPROVED

MMDadoda
 10/1/2015 2:24:37 PM

Quant Time: Oct 01 06:21:30 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 01 05:22:05 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	119221	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	453081	20.00	ng	0.00
38) Acenaphthene-d10	9.94	164	185752	20.00	ng	-0.01
63) Phenanthrene-d10	11.43	188	290845	20.00	ng	-0.01
75) Chrysene-d12	14.07	240	306368	20.00	ng	-0.01
86) Perylene-d12	15.53	264	309644	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	801649	122.06	ng	0.01
7) Phenol-d6	6.54	99	981641	118.78	ng	0.00
23) Nitrobenzene-d5	7.47	82	659779	97.07	ng	0.00
41) 2,4,6-Tribromophenol	10.74	330	230920	122.75	ng	0.00
44) 2-Fluorobiphenyl	9.27	172	1077393	100.61	ng	0.00
78) Terphenyl-d14	13.02	244	853597	92.10	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.67	163	322774	25.17	ng	Qvalue # 97
74) Fluoranthene	12.64	202	108911	6.99	ng	95
77) Pyrene	12.87	202	113384	6.19	ng	97
80) Benzo(a)anthracene	14.07	228	95446	5.79	ng	# 91
82) Chrysene	14.10	228	79706m	4.57	ng	
85) Indeno(1,2,3-cd)pyrene	16.97	276	46628	3.61	ng	# 85
87) Benzo(b)fluoranthene	15.11	252	120811m	6.83	ng	
88) Benzo(k)fluoranthene	15.12	252	64876m	4.00	ng	
89) Benzo(a)pyrene	15.48	252	81986	5.35	ng	# 85
91) Benzo(g,h,i)perylene	17.41	276	40201	3.05	ng	# 74

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

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