

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032116\
 Data File : BF085661.D
 Acq On : 22 Mar 2016 11:44
 Operator : UM/SJ
 Sample : H1799-10 10X
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AOU2-SB37A(COMP7-21)

Manual Integrations
 APPROVED

Sohil
 3/22/2016 5:34:54 PM

Quant Time: Mar 22 13:19:56 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 18 23:31:39 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	114385	20.00	ng	-0.02
21) Naphthalene-d8	8.13	136	480601	20.00	ng	-0.01
38) Acenaphthene-d10	9.88	164	192647	20.00	ng	-0.02
63) Phenanthrene-d10	11.36	188	330492	20.00	ng	-0.02
75) Chrysene-d12	13.99	240	226755	20.00	ng	-0.02
86) Perylene-d12	15.42	264	221367	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	80537	11.83	ng	-0.02
7) Phenol-d6	6.47	99	103502	11.85	ng	-0.02
23) Nitrobenzene-d5	7.40	82	59546	7.19	ng	-0.03
41) 2,4,6-Tribromophenol	10.66	330	17483	9.26	ng	-0.02
44) 2-Fluorobiphenyl	9.20	172	130637	8.26	ng	-0.02
78) Terphenyl-d14	12.94	244	81949	8.70	ng	-0.02
Target Compounds						Qvalue
31) Naphthalene	8.15	128	49124	2.03	ng	96
57) Fluorene	10.42	166	63024	5.04	ng	99
70) Phenanthrene	11.38	178	137152	7.87	ng	99
74) Fluoranthene	12.57	202	90061	4.94	ng	92
77) Pyrene	12.80	202	85289	5.24	ng	97
80) Benzo(a)anthracene	13.98	228	48315	3.67	ng	97
82) Chrysene	14.01	228	35704m	2.82	ng	
85) Indeno(1,2,3-cd)pyrene	16.80	276	22071	2.09	ng	92
87) Benzo(b)fluoranthene	15.01	252	59166m	4.10	ng	
89) Benzo(a)pyrene	15.36	252	38501	3.14	ng	# 95

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

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