

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062716\
 Data File : BF088479.D
 Acq On : 28 Jun 2016 6:04
 Operator : UM/SJ
 Sample : H3624-05DL 2X
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SC-STA-1481(0-4)DL

Quant Time: Jun 28 07:34:34 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 28 05:25:55 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.54	152	83349m	20.00	ng	-0.01
21) Naphthalene-d8	7.83	136	329782	20.00	ng	-0.01
38) Acenaphthene-d10	9.58	164	135190	20.00	ng	-0.01
63) Phenanthrene-d10	11.05	188	200467	20.00	ng	-0.01
75) Chrysene-d12	13.69	240	158879	20.00	ng	0.00
86) Perylene-d12	15.05	264	172607	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.12	112	235600	48.32	ng	-0.01
7) Phenol-d6	6.20	99	328378	55.57	ng	-0.02
23) Nitrobenzene-d5	7.12	82	177070	31.35	ng	-0.01
41) 2,4,6-Tribromophenol	10.36	330	44256	31.00	ng	-0.01
44) 2-Fluorobiphenyl	8.90	172	358962	38.70	ng	-0.02
78) Terphenyl-d14	12.64	244	193509	27.72	ng	0.00
Target Compounds						Qvalue
48) Acenaphthylene	9.44	152	144085	10.35	ng	98
49) Dimethylphthalate	9.31	163	48567	4.96	ng	98
70) Phenanthrene	11.07	178	169942	16.16	ng	99
71) Anthracene	11.13	178	159381	15.02	ng	99
72) Carbazole	11.30	167	30437	3.07	ng	94
74) Fluoranthene	12.26	202	584877	52.69	ng	99
77) Pyrene	12.49	202	673838	57.15	ng	100
80) Benzo(a)anthracene	13.68	228	536143	59.22	ng	94
82) Chrysene	13.71	228	345935m	40.61	ng	
85) Indeno(1,2,3-cd)pyrene	16.29	276	196321	24.81	ng	# 100
87) Benzo(b)fluoranthene	14.69	252	550009m	45.72	ng	
88) Benzo(k)fluoranthene	14.70	252	291453m	32.12	ng	
89) Benzo(a)pyrene	14.99	252	385277	39.58	ng	99
90) Dibenzo(a,h)anthracene	16.27	278	51084m	5.65	ng	
91) Benzo(a,h,i)perylene	16.65	276	158273	17.02	ng	98

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

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