

Data Path : U:\HPCHEM1\BNA F\DATA\BF082217\
 Data File : BF097965.D
 Acq On : 22 Aug 2017 20:38
 Operator : SJ/JU
 Sample : I4889-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SASP2P-TP-106-E(0-5)COMP

Manual Integrations
 APPROVED

Sohil
 8/23/2017 3:03:47 PM

Quant Time: Aug 23 01:44:49 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 21 15:59:04 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	143080	20.00	ng	0.00
21) Naphthalene-d8	8.03	136	562769	20.00	ng	0.00
38) Acenaphthene-d10	9.78	164	241796	20.00	ng	0.00
63) Phenanthrene-d10	11.26	188	405147	20.00	ng	0.00
75) Chrysene-d12	13.90	240	287577	20.00	ng	0.00
86) Perylene-d12	15.32	264	270125	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	562240	59.77	ng	0.00
7) Phenol-d6	6.38	99	750373	58.71	ng	0.00
23) Nitrobenzene-d5	7.31	82	587093	56.67	ng	-0.01
41) 2,4,6-Tribromophenol	10.57	330	116813	55.68	ng	0.00
44) 2-Fluorobiphenyl	9.11	172	862357	52.40	ng	0.00
78) Terphenyl-d14	12.85	244	542731	39.36	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.50	163	72110	4.080	ng	97
74) Fluoranthene	12.47	202	67592	3.000	ng	98
77) Pyrene	12.70	202	69889	2.971	ng	95
80) Benzo(a)anthracene	13.89	228	42741	2.480	ng	93
82) Chrysene	13.92	228	42981m	2.507	ng	
85) Indeno(1,2,3-cd)pyrene	16.69	276	34016	2.697	ng	# 90
87) Benzo(b)fluoranthene	14.91	252	75576m	4.439	ng	
89) Benzo(a)pyrene	15.26	252	63574	4.218	ng	97
91) Benzo(g,h,i)perylene	17.10	276	34668	2.708	ng	92

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

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