

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080717\
 Data File : BF097442.D
 Acq On : 7 Aug 2017 21:20
 Operator : SJ/JU
 Sample : I4623-05
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-WE

Manual Integrations
 APPROVED

Sohil
 8/9/2017 7:56:03 PM

Quant Time: Aug 08 05:47:02 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 07 16:02:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.41	152	89030	20.00	ng	0.00
21) Naphthalene-d8	7.69	136	410322	20.00	ng	0.00
38) Acenaphthene-d10	9.43	164	146575	20.00	ng	0.00
63) Phenanthrene-d10	10.91	188	205864	20.00	ng	0.00
75) Chrysene-d12	13.53	240	191356	20.00	ng	0.00
86) Perylene-d12	14.87	264	158956	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.00	112	579258	98.08	ng	0.00
7) Phenol-d6	6.08	99	657876	97.58	ng	0.00
23) Nitrobenzene-d5	6.98	82	456001	65.19	ng	-0.01
41) 2,4,6-Tribromophenol	10.23	330	106594	92.04	ng	0.00
44) 2-Fluorobiphenyl	8.77	172	680527	78.79	ng	0.00
78) Terphenyl-d14	12.49	244	403002	42.94	ng	0.00
Target Compounds						
10) Phenol	6.09	94	20549	2.569	ng	Qvalue # 55
49) Dimethylphthalate	9.17	163	160914	14.457	ng	98
51) Acenaphthene	9.47	154	19674	2.362	ng	96
57) Fluorene	9.98	166	16970	2.035	ng	92
70) Phenanthrene	10.93	178	184624	16.738	ng	99
71) Anthracene	10.99	178	40555	3.590	ng	99
74) Fluoranthene	12.11	202	245006	22.673	ng	96
77) Pyrene	12.34	202	195094	12.454	ng	99
80) Benzo(a)anthracene	13.52	228	111887	9.037	ng	99
82) Chrysene	13.56	228	94486	7.815	ng	97
85) Indeno(1,2,3-cd)pyrene	16.06	276	34711	3.348	ng	98
87) Benzo(b)fluoranthene	14.53	252	117683m	12.316	ng	
88) Benzo(k)fluoranthene	14.54	252	38182m	4.193	ng	
89) Benzo(a)pyrene	14.82	252	79359	9.075	ng	99
91) Benzo(a,h,i)perylene	16.42	276	29837	3.967	ng	93

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

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