

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080317\
 Data File : BF097344.D
 Acq On : 4 Aug 2017 4:04
 Operator : SJ/JU
 Sample : I4522-02
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-3D68

Manual Integrations
 APPROVED

Sohil
 8/4/2017 5:31:06 PM

Quant Time: Aug 04 11:10:21 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 04 01:17:22 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.45	152	95212	20.00	ng	-0.05
21) Naphthalene-d8	7.73	136	336066	20.00	ng	-0.06
38) Acenaphthene-d10	9.47	164	143433	20.00	ng	-0.06
63) Phenanthrene-d10	10.94	188	202529	20.00	ng	-0.06
75) Chrysene-d12	13.58	240	117688	20.00	ng	-0.05
86) Perylene-d12	14.93	264	92547	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.05	112	600520	95.08	ng	-0.05
7) Phenol-d6	6.12	99	692336	96.02	ng	-0.05
23) Nitrobenzene-d5	7.02	82	376390	65.70	ng	-0.06
41) 2,4,6-Tribromophenol	10.26	330	94262	83.18	ng	-0.06
44) 2-Fluorobiphenyl	8.80	172	447902	53.00	ng	-0.06
78) Terphenyl-d14	12.53	244	274359	47.53	ng	-0.05
Target Compounds						
10) Phenol	6.14	94	55144	6.448	ng	Qvalue # 66
31) Naphthalene	7.75	128	227810	14.380	ng	100
37) 2-Methylnaphthalene	8.44	142	51182	5.103	ng	98
48) Acenaphthylene	9.33	152	47119	3.405	ng	99
49) Dimethylphthalate	9.20	163	91711	8.420	ng	100
51) Acenaphthene	9.50	154	95148	11.673	ng	99
54) Dibenzofuran	9.67	168	93636	8.624	ng	98
57) Fluorene	10.02	166	109808	13.457	ng	99
70) Phenanthrene	10.97	178	819482	75.517	ng	99
71) Anthracene	11.02	178	260733	23.459	ng	98
72) Carbazole	11.18	167	158761	14.524	ng	99
74) Fluoranthene	12.16	202	964557	90.732	ng	98
77) Pvrene	12.39	202	778464	80.798	ng	99
80) Benzo(a)anthracene	13.57	228	376985	49.506	ng	98
82) Chrysene	13.61	228	322188	43.330	ng	96
85) Indeno(1,2,3-cd)pyrene	16.13	276	146050	22.906	ng	97
87) Benzo(b)fluoranthene	14.57	252	330377m	59.386	ng	
88) Benzo(k)fluoranthene	14.59	252	133232m	25.132	ng	
89) Benzo(a)pvrene	14.87	252	245899	48.295	ng	# 95
90) Dibenzo(a,h)anthracene	16.13	278	41332m	9.899	ng	
91) Benzo(q,h,i)perylene	16.50	276	139381	31.833	ng	97

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

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