

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF021125\
 Data File : BF141568.D
 Acq On : 11 Feb 2025 18:01
 Operator : RC/JU
 Sample : Q1285-01DL 100X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 72-11977DL

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 02/12/2025
 Supervised By :mohammad ahmed 02/12/2025

Quant Time: Feb 11 18:22:37 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF020625.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 06 16:58:59 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.786	152	73262	20.000	ng	0.00
21) Naphthalene-d8	8.069	136	283534	20.000	ng	0.00
39) Acenaphthene-d10	9.822	164	144692	20.000	ng	-0.01
64) Phenanthrene-d10	11.304	188	212403	20.000	ng	-0.01
76) Chrysene-d12	13.945	240	171943	20.000	ng	-0.02
86) Perylene-d12	15.404	264	199200	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	3069	0.657	ng	0.01
7) Phenol-d6	6.439	99	4248	0.719	ng	0.00
23) Nitrobenzene-d5	7.351	82	3235	0.595	ng	-0.01
42) 2,4,6-Tribromophenol	10.616	330	739	0.508	ng	-0.01
45) 2-Fluorobiphenyl	9.139	172	5631	0.600	ng	-0.01
79) Terphenyl-d14	12.886	244	5948	0.584	ng	-0.02
Target Compounds						Qvalue
31) Naphthalene	8.092	128	624724	42.328	ng	100
37) 2-Methylnaphthalene	8.780	142	276935	28.835	ng	100
38) 1-Methylnaphthalene	8.880	142	225723	24.266	ng	97
46) 1,1'-Biphenyl	9.245	154	90597	8.076	ng	98
49) Acenaphthylene	9.680	152	61292	4.968	ng	# 95
52) Acenaphthene	9.851	154	170769	20.588	ng	99
58) Fluorene	10.369	166	104691	11.121	ng	97
71) Phenanthrene	11.333	178	472019	40.372	ng	99
72) Anthracene	11.380	178	146047	12.426	ng	99
75) Fluoranthene	12.516	202	139667	11.267	ng	100
78) Pyrene	12.745	202	274343	18.743	ng	99
81) Benzo(a)anthracene	13.939	228	90910m	7.954	ng	
83) Chrysene	13.974	228	65279	6.186	ng	100
87) Indeno(1,2,3-cd)pyrene	16.839	276	26517	2.154	ng	95
88) Benzo(b)fluoranthene	14.992	252	61536m	4.601	ng	
90) Benzo(a)pyrene	15.345	252	82036	7.580	ng	97
92) Benzo(g,h,i)perylene	17.268	276	28978	2.777	ng	98

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

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