

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061824\
 Data File : BF138222.D
 Acq On : 18 Jun 2024 14:02
 Operator : CG\JU
 Sample : P2924-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DAMAGED-CONC

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 06/19/2024
 Supervised By :mohammad ahmed 06/21/2024

Quant Time: Jun 18 14:26:12 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 13 01:52:47 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.892	152	292730	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	990558	20.000	ng	0.00
39) Acenaphthene-d10	10.004	164	316383	20.000	ng	# 0.08
64) Phenanthrene-d10	11.574	188	111639	20.000	ng	# 0.16
76) Chrysene-d12	14.115	240	422310	20.000	ng	# 0.06
86) Perylene-d12	15.533	264	294812	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	987479	56.430	ng	0.00
7) Phenol-d6	6.516	99	1565557	70.554	ng	0.00
23) Nitrobenzene-d5	7.451	82	1168106	68.651	ng	0.00
42) 2,4,6-Tribromophenol	10.822	330	66778	21.202	ng	0.11
45) 2-Fluorobiphenyl	9.275	172	2105930	105.845	ng	0.02
79) Terphenyl-d14	13.127	244	474235	17.786	ng	0.13
Target Compounds						
						Qvalue
31) Naphthalene	8.192	128	699261	14.276	ng	99
37) 2-Methylnaphthalene	8.898	142	111816	3.478	ng	100
38) 1-Methylnaphthalene	9.016	142	105887	3.361	ng	96
46) 1,1'-Biphenyl	9.398	154	195818	8.435	ng	78
88) Benzo(b)fluoranthene	15.110	252	54781m	3.151	ng	

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

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