

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010323\
 Data File : BG056195.D
 Acq On : 3 Jan 2023 21:07
 Operator : CG/JU
 Sample : N6244-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PE-13

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 01/04/2023
 Supervised By :Yogesh Patel 01/05/2023

Quant Time: Jan 04 02:16:11 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 04 02:08:15 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.347	152	62506	20.000	ng	0.00
21) Naphthalene-d8	11.173	136	232012	20.000	ng	0.00
39) Acenaphthene-d10	14.945	164	174186	20.000	ng	0.00
64) Phenanthrene-d10	17.683	188	402541	20.000	ng	0.00
76) Chrysene-d12	21.983	240	366650	20.000	ng	0.00
86) Perylene-d12	25.450	264	394707	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.861	112	282396	80.667	ng	0.01
7) Phenol-d6	7.477	99	399957	74.394	ng	0.00
23) Nitrobenzene-d5	9.510	82	224053	49.397	ng	0.00
42) 2,4,6-Tribromophenol	16.425	330	153877	69.753	ng	0.00
45) 2-Fluorobiphenyl	13.576	172	660674	52.376	ng	0.00
79) Terphenyl-d14	20.256	244	1005947	49.140	ng	0.00
Target Compounds						Qvalue
71) Phenanthrene	17.724	178	112474	5.392	ng	98
75) Fluoranthene	19.716	202	279830	10.284	ng	99
78) Pyrene	20.074	202	247554	9.579	ng	99
81) Benzo(a)anthracene	21.966	228	157887	6.131	ng	98
83) Chrysene	22.036	228	133403	5.530	ng	97
87) Indeno(1,2,3-cd)pyrene	29.428	276	90646	3.158	ng	# 91
88) Benzo(b)fluoranthene	24.340	252	184568	7.360	ng	95
89) Benzo(k)fluoranthene	24.404	252	57405m	2.299	ng	
90) Benzo(a)pyrene	25.280	252	134672	5.506	ng	99
92) Benzo(g,h,i)perylene	30.679	276	86401	3.702	ng	97

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

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