

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122021\
 Data File : BF126275.D
 Acq On : 20 Dec 2021 13:31
 Operator : JU/CG
 Sample : M5124-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 F2337-STOPJOINT-TP

Quant Time: Dec 20 14:29:34 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF121521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 16 12:22:48 2021
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Christian Giraldo 12/21/2021
 Supervised By :Jagrut Upadhyay 12/21/2021

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.816	152	174925	20.000	ng	0.00
21) Naphthalene-d8	8.098	136	728013	20.000	ng	# 0.00
39) Acenaphthene-d10	9.851	164	300309	20.000	ng	0.00
64) Phenanthrene-d10	11.333	188	411817	20.000	ng	# 0.00
76) Chrysene-d12	13.968	240	274046	20.000	ng	# 0.00
86) Perylene-d12	15.415	264	313163	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	1312833	110.524	ng	0.01
7) Phenol-d6	6.445	99	1611070	106.818	ng	0.00
23) Nitrobenzene-d5	7.375	82	1053871	78.401	ng	0.00
42) 2,4,6-Tribromophenol	10.639	330	236741	98.094	ng	0.00
45) 2-Fluorobiphenyl	9.175	172	1374730	80.288	ng	0.00
79) Terphenyl-d14	12.921	244	897042	56.894	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.451	94	36657	2.165	ng	95
75) Fluoranthene	12.545	202	50044	2.613	ng	94
78) Pyrene	12.774	202	55152	2.191	ng	98
88) Benzo(b)fluoranthene	14.998	252	48453m	2.575	ng	
90) Benzo(a)pyrene	15.351	252	37037	2.368	ng	# 94

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

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