

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062422\
 Data File : BF128989.D
 Acq On : 24 Jun 2022 15:02
 Operator : CG\JU
 Sample : N3424-03RE
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OH-CON-0621RE

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 06/27/2022
 Supervised By :Jagrut Upadhyay 06/27/2022

Quant Time: Jun 25 06:13:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 24 00:50:48 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.722	152	113271	20.000	ng	0.00
21) Naphthalene-d8	8.004	136	431286	20.000	ng	0.00
39) Acenaphthene-d10	9.763	164	222790	20.000	ng	0.00
64) Phenanthrene-d10	11.251	188	321899	20.000	ng	0.00
76) Chrysene-d12	13.898	240	183943	20.000	ng	# 0.00
86) Perylene-d12	15.339	264	159471	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.387	112	1392	0.196	ng	0.03
7) Phenol-d6	6.392	99	17270	1.990	ng	0.00
23) Nitrobenzene-d5	7.287	82	219541	23.700	ng	-0.02
42) 2,4,6-Tribromophenol	0.000	330	0	0.000	ng	
45) 2-Fluorobiphenyl	9.081	172	422822	26.585	ng	-0.01
79) Terphenyl-d14	12.833	244	310902	29.370	ng	0.00
Target Compounds						Qvalue
78) Pyrene	12.692	202	46767	3.417	ng	98
88) Benzo(b)fluoranthene	14.927	252	33020m	3.240	ng	
90) Benzo(a)pyrene	15.280	252	22234	2.776	ng	# 87

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

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