

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF041522\
 Data File : BF127800.D
 Acq On : 15 Apr 2022 12:45
 Operator : CG\JU
 Sample : N2389-07RE
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 33637-8RE

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 04/18/2022
 Supervised By : Jagrut Upadhyay 04/18/2022

Quant Time: Apr 18 01:23:16 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF040622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 13 16:40:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.951	152	131076	20.000	ng	-0.01
21) Naphthalene-d8	8.233	136	436090	20.000	ng	#-0.01
39) Acenaphthene-d10	9.998	164	183452	20.000	ng	0.00
64) Phenanthrene-d10	11.486	188	303888	20.000	ng	0.00
76) Chrysene-d12	14.133	240	316457	20.000	ng	-0.01
86) Perylene-d12	15.651	264	351057	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.587	112	784101	97.202	ng	0.00
7) Phenol-d6	6.587	99	856392	81.152	ng	0.00
23) Nitrobenzene-d5	7.516	82	723781	73.992	ng	-0.01
42) 2,4,6-Tribromophenol	10.798	330	216982	111.147	ng	0.00
45) 2-Fluorobiphenyl	9.310	172	1182799	91.210	ng	-0.01
79) Terphenyl-d14	13.074	244	1139363	59.428	ng	0.00
Target Compounds						Qvalue
9) Benzaldehyde	6.504	77	78134	8.740	ng	93
10) Phenol	6.598	94	140858	13.349	ng	100
15) Benzyl Alcohol	7.098	79	87137	9.445	ng	97
17) 2-Methylphenol	7.198	107	43586	6.123	ng	97
22) Acetophenone	7.357	105	261329	22.301	ng	# 81
27) 2,4-Dimethylphenol	7.886	122	18860	2.827	ng	83
31) Naphthalene	8.257	128	1380471	60.411	ng	98
32) Benzoic acid	8.175	122	2119984m	437.968	ng	
37) 2-Methylnaphthalene	8.945	142	37429	2.449	ng	95

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

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