

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071222\
 Data File : BF129194.D
 Acq On : 12 Jul 2022 17:45
 Operator : CG\JU
 Sample : N3663-02
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VB16181

Manual Integrations
 APPROVED

Reviewed By :Christian Giraldo 07/13/2022
 Supervised By :Jagrut Upadhyay 07/13/2022

Quant Time: Jul 12 18:39:26 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 12 14:13:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.934	152	338422	20.000	ng	0.00
21) Naphthalene-d8	8.216	136	1319437	20.000	ng	0.00
39) Acenaphthene-d10	9.974	164	629727	20.000	ng	0.00
64) Phenanthrene-d10	11.463	188	1003580	20.000	ng	# 0.00
76) Chrysene-d12	14.110	240	568858	20.000	ng	-0.01
86) Perylene-d12	15.621	264	465379	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.575	112	2272780	113.540	ng	0.01
7) Phenol-d6	6.563	99	2553772	102.722	ng	0.00
23) Nitrobenzene-d5	7.498	82	2052954	92.806	ng	0.00
42) 2,4,6-Tribromophenol	10.769	330	826160	120.654	ng	0.00
45) 2-Fluorobiphenyl	9.292	172	3297339	83.767	ng	0.00
79) Terphenyl-d14	13.057	244	2826562	103.184	ng	0.00
Target Compounds						
20) 3+4-Methylphenols	7.328	107	226471	10.406	ng	# 70
32) Benzoic acid	7.933	122	57321m	3.978	ng	

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

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