

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061722\
 Data File : BF128856.D
 Acq On : 17 Jun 2022 16:09
 Operator : CG\JU
 Sample : N3373-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SC-1W

Quant Time: Jun 18 05:06:42 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 14 14:53:53 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.781	152	84463	20.000	ng	0.00
21) Naphthalene-d8	8.063	136	290617	20.000	ng	0.00
39) Acenaphthene-d10	9.822	164	148696	20.000	ng	0.00
64) Phenanthrene-d10	11.310	188	235645	20.000	ng	# 0.00
76) Chrysene-d12	13.957	240	170502	20.000	ng	# 0.00
86) Perylene-d12	15.410	264	167450	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	463659	87.438	ng	0.00
7) Phenol-d6	6.440	99	649141	100.287	ng	0.00
23) Nitrobenzene-d5	7.345	82	459135	73.556	ng	0.00
42) 2,4,6-Tribromophenol	10.616	330	32536	18.519	ng	0.00
45) 2-Fluorobiphenyl	9.145	172	798265	75.201	ng	0.00
79) Terphenyl-d14	12.898	244	618326	63.015	ng	0.00
Target Compounds						
						Qvalue
55) Dibenzofuran	10.022	168	37443	3.070	ng	93
71) Phenanthrene	11.339	178	994640	84.153	ng	98
73) Carbazole	11.545	167	318944	29.353	ng	99
75) Fluoranthene	12.551	202	3634819	294.201	ng	98
78) Pyrene	12.774	202	1981839	156.231	ng	99
81) Benzo(a)anthracene	13.945	228	122399	11.517	ng	97
83) Chrysene	13.980	228	122646	12.104	ng	99

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

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