

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062022\
 Data File : BF128879.D
 Acq On : 20 Jun 2022 16:24
 Operator : CG\JU
 Sample : N3373-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SC-1N

Quant Time: Jun 21 01:15:59 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 21 00:40:09 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.769	152	110080	20.000	ng	0.00
21) Naphthalene-d8	8.057	136	417122	20.000	ng	# 0.00
39) Acenaphthene-d10	9.810	164	235798	20.000	ng	0.00
64) Phenanthrene-d10	11.304	188	412538	20.000	ng	# 0.00
76) Chrysene-d12	13.945	240	221163	20.000	ng	# 0.00
86) Perylene-d12	15.392	264	198506	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.410	112	565487	81.825	ng	0.00
7) Phenol-d6	6.428	99	813962	96.486	ng	-0.01
23) Nitrobenzene-d5	7.340	82	584879	65.283	ng	-0.01
42) 2,4,6-Tribromophenol	10.604	330	66661	23.927	ng	0.00
45) 2-Fluorobiphenyl	9.133	172	1152786	68.483	ng	0.00
79) Terphenyl-d14	12.892	244	1104363	86.768	ng	0.00
Target Compounds						
						Qvalue
55) Dibenzofuran	10.016	168	62250	3.218	ng	96
71) Phenanthrene	11.333	178	2037254	98.456	ng	100
73) Carbazole	11.533	167	287496	15.114	ng	98
75) Fluoranthene	12.539	202	3448908	159.455	ng	98
78) Pyrene	12.757	202	1708859	103.854	ng	99
81) Benzo(a)anthracene	13.933	228	46744	3.391	ng	98
83) Chrysene	13.968	228	113322	8.622	ng	97

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

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