

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062222\
 Data File : BF128936.D
 Acq On : 22 Jun 2022 17:25
 Operator : CG\JU
 Sample : N3373-07DL 10X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SC-6EDL

Quant Time: Jun 23 01:30:37 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 22 01:02:29 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.745	152	127472	20.000	ng	-0.01
21) Naphthalene-d8	8.028	136	487059	20.000	ng	-0.01
39) Acenaphthene-d10	9.786	164	283089	20.000	ng	0.00
64) Phenanthrene-d10	11.274	188	508464	20.000	ng	0.00
76) Chrysene-d12	13.915	240	389056	20.000	ng	#-0.01
86) Perylene-d12	15.357	264	306996	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.387	112	55425	6.926	ng	0.00
7) Phenol-d6	6.404	99	82036	8.398	ng	-0.01
23) Nitrobenzene-d5	7.310	82	57188	5.467	ng	-0.02
42) 2,4,6-Tribromophenol	10.575	330	8678	2.594	ng	-0.01
45) 2-Fluorobiphenyl	9.104	172	122994	6.086	ng	-0.01
79) Terphenyl-d14	12.857	244	138936	6.205	ng	-0.01
Target Compounds						
						Qvalue
71) Phenanthrene	11.298	178	898631	35.236	ng	99
73) Carbazole	11.510	167	78850	3.363	ng	97
75) Fluoranthene	12.492	202	1117848	41.932	ng	97
78) Pyrene	12.721	202	748038	25.843	ng	99

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

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