

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111022\
 Data File : BM037455.D
 Acq On : 10 Nov 2022 17:22
 Operator : CG/JU
 Sample : N5495-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SB-29A

Quant Time: Nov 11 00:18:14 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 09 02:08:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	215696	20.000	ng	-0.01
21) Naphthalene-d8	10.610	136	852909	20.000	ng	-0.02
39) Acenaphthene-d10	14.451	164	503422	20.000	ng	-0.01
64) Phenanthrene-d10	17.192	188	1050460	20.000	ng	-0.01
76) Chrysene-d12	21.374	240	1083147	20.000	ng	0.00
86) Perylene-d12	23.703	264	1136260	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.387	112	1309000	104.846	ng	0.00
7) Phenol-d6	6.992	99	1648385	97.279	ng	-0.01
23) Nitrobenzene-d5	8.980	82	1001442	59.239	ng	-0.02
42) 2,4,6-Tribromophenol	15.939	330	722903	121.851	ng	-0.02
45) 2-Fluorobiphenyl	13.074	172	2235674	59.007	ng	-0.02
79) Terphenyl-d14	19.815	244	3757285	62.206	ng	-0.01
Target Compounds						Qvalue
32) Benzoic acid	9.945	122	418196	41.465	ng	98
70) Pentachlorophenol	16.851	266	25348	3.054	ng	97
75) Fluoranthene	19.250	202	140456	2.039	ng	99
78) Pyrene	19.609	202	188856	2.267	ng	100
80) Butylbenzylphthalate	20.509	149	96993	3.261	ng	99
88) Benzo(b)fluoranthene	22.991	252	164525	2.086	ng	# 90

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

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