

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102323\
 Data File : BM042410.D
 Acq On : 23 Oct 2023 19:59
 Operator : MA/JU
 Sample : 04945-01
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PAT-2-101823

Quant Time: Oct 23 22:33:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 21 02:32:53 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.993	152	134984	20.000	ng	0.00
21) Naphthalene-d8	10.816	136	522751	20.000	ng	0.00
39) Acenaphthene-d10	14.639	164	264882	20.000	ng	0.00
64) Phenanthrene-d10	17.392	188	566236	20.000	ng	0.00
76) Chrysene-d12	21.574	240	591544	20.000	ng	0.00
86) Perylene-d12	24.039	264	676141	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	906129	94.874	ng	0.00
7) Phenol-d6	7.146	99	1179619	92.162	ng	0.00
23) Nitrobenzene-d5	9.192	82	724529	61.288	ng	0.00
42) 2,4,6-Tribromophenol	16.133	330	261736	92.355	ng	0.00
45) 2-Fluorobiphenyl	13.263	172	1183459	60.165	ng	0.00
79) Terphenyl-d14	20.004	244	1938281	58.111	ng	0.00
Target Compounds						Qvalue
74) Di-n-butylphthalate	18.333	149	5253	2.954	ng	# 91
75) Fluoranthene	19.445	202	79799	2.352	ng	97
78) Pyrene	19.809	202	90956	2.200	ng	98
84) Bis(2-ethylhexyl)phtha...	21.439	149	7203	3.620	ng	# 100
88) Benzo(b)fluoranthene	23.274	252	139206	3.420	ng	# 96
90) Benzo(a)pyrene	23.927	252	101801	2.599	ng	# 97
92) Benzo(g,h,i)perylene	27.427	276	94473	2.368	ng	97

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

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