

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110123\
 Data File : BM042540.D
 Acq On : 01 Nov 2023 15:24
 Operator : MA/JU
 Sample : 05138-03RE
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 348RE

Quant Time: Nov 01 16:38:00 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM103023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 31 02:55:18 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	68377	20.000	ng	0.00
21) Naphthalene-d8	10.739	136	254497	20.000	ng	-0.01
39) Acenaphthene-d10	14.568	164	107924	20.000	ng	-0.01
64) Phenanthrene-d10	17.321	188	349496	20.000	ng	0.00
76) Chrysene-d12	21.503	240	238602	20.000	ng	0.00
86) Perylene-d12	23.921	264	22003	20.000	ng	#-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	630981	148.928	ng	0.00
7) Phenol-d6	7.092	99	651492	112.121	ng	0.00
23) Nitrobenzene-d5	9.122	82	610836	111.887	ng	-0.01
42) 2,4,6-Tribromophenol	16.068	330	396738	281.566	ng	0.00
45) 2-Fluorobiphenyl	13.192	172	1271239	158.498	ng	0.00
79) Terphenyl-d14	19.939	244	2521961	179.221	ng	0.00
Target Compounds						
60) Diethylphthalate	15.380	149	31326	3.833	ng	98
74) Di-n-butylphthalate	18.262	149	101327	4.903	ng	98
84) Bis(2-ethylhexyl)phtha...	21.368	149	63899	6.042	ng	99
88) Benzo(b)fluoranthene	23.174	252	13640	11.127	ng	# 74
89) Benzo(k)fluoranthene	23.221	252	10580	7.745	ng	# 69

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

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