

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060820\
 Data File : BM026290.D
 Acq On : 08 Jun 2020 17:08
 Operator : JU/CG
 Sample : L2904-07
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 VE-3-1

Manual Integrations
 APPROVED

mohammad
 6/9/2020 12:21:43 PM

Quant Time: Jun 08 17:45:47 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM060520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 08 15:00:18 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.45	152	206483	20.00	ng	0.00
21) Naphthalene-d8	10.20	136	789843	20.00	ng	0.00
39) Acenaphthene-d10	14.10	164	425222	20.00	ng	0.00
64) Phenanthrene-d10	16.85	188	670951	20.00	ng	0.00
76) Chrysene-d12	21.06	240	657919	20.00	ng	0.00
86) Perylene-d12	23.16	264	706919	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.08	112	574171	49.16	ng	0.00
7) Phenol-d6	6.65	99	506192	29.93	ng	0.00
23) Nitrobenzene-d5	8.59	82	1269260	78.90	ng	0.00
42) 2,4,6-Tribromophenol	15.60	330	547027	106.20	ng	0.01
45) 2-Fluorobiphenyl	12.71	172	2445907	87.31	ng	0.00
79) Terphenyl-d14	19.50	244	2736105	80.73	ng	0.00
Target Compounds						
20) 3+4-Methylphenols	8.23	107	214591	11.961	ng	94
27) 2,4-Dimethylphenol	9.45	122	72021m	6.845	ng	
50) Dimethylphthalate	13.57	163	285156	9.168	ng	# 95
52) Acenaphthene	14.16	154	80576	3.466	ng	98
58) Fluorene	15.15	166	78002	2.563	ng	# 86

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

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